

SEARCH REQUEST FORM

Scientific and Technical Information Center

Requester's Full Name: Hong Liu Examiner #: _____ Date: 8/2/02
 Art Unit: 1624 Phone Number 30 6-5814 Serial Number: 09/864,905
 Mail Box and Bldg/Room Location: 4C01 Results Format Preferred (circle) PAPER ~~DISK~~ ~~E-MAIL~~
4C12

If more than one search is submitted, please prioritize searches in order of need.

Please provide a detailed statement of the search topic, and describe as specifically as possible the subject matter to be searched. Include the elected species or structures, keywords, synonyms, acronyms, and registry numbers, and combine with the concept or utility of the invention. Define any terms that may have a special meaning. Give examples or relevant citations, authors, etc, if known. Please attach a copy of the cover sheet, pertinent claims, and abstract.

Title of Invention: Pharm. compositions & methods for use
 Inventors (please provide full names): Miller, C Dull, G Miao, L Lynn, D
Schmitt, J Clark, T

Earliest Priority Filing Date: _____

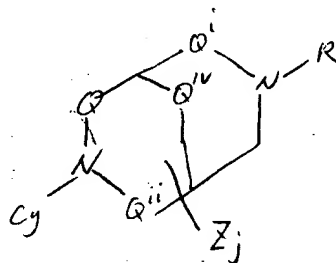
For Sequence Searches Only Please include all pertinent information (parent, child, divisional, or issued patent numbers) along with the appropriate serial number.

Barb please!

RECEIVED

AUG 21 2002

STIC



Point of Contact:
 Barb O'Brien
 Technical Information Specialist
 STIC CM1 6A05 308-4291

Applicants elected the species of 3-(5-phenoxyl-3-pyridyl)-
 3,7-diazabicyclo[3.3.1]nonane. Please based the structure search
 on this core.

STAFF USE ONLY

Searcher: BOB
 Searcher Phone #: _____
 Searcher Location: _____
 Date Searcher Picked Up: _____
 Date Completed: 8-30-02
 Searcher Prep & Review Time: 40
 Clerical Prep Time: _____
 Online Time: 35

Type of Search

NA Sequence (#) _____
 AA Sequence (#) _____
 Structure (#) 3
 Bibliographic _____
 Litigation _____
 Fulltext _____
 Patent Family _____
 Other _____

Vendors and cost where applicable

STN 242
 Dialog _____
 Questel/Orbit _____
 Dr.Link _____
 Lexis/Nexis _____
 Sequence Systems _____
 WWW/Internet _____
 Other (specify) Chem Draw

Welcome to STN International! Enter x:x

LOGINID:sssptal611hxl

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

species for w097/40.49

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Apr 08 "Ask CAS" for self-help around the clock
NEWS 3 Apr 09 BEILSTEIN: Reload and Implementation of a New Subject Area
NEWS 4 Apr 09 ZDB will be removed from STN
NEWS 5 Apr 19 US Patent Applications available in IFICDB, IFIPAT, and
IFIUDB
NEWS 6 Apr 22 Records from IP.com available in CAPLUS, HCAPLUS, and
ZCAPLUS
NEWS 7 Apr 22 BIOSIS Gene Names now available in TOXCENTER
NEWS 8 Apr 22 Federal Research in Progress (FEDRIP) now available
NEWS 9 Jun 03 New e-mail delivery for search results now available
NEWS 10 Jun 10 MEDLINE Reload
NEWS 11 Jun 10 PCTFULL has been reloaded
NEWS 12 Jul 02 FOREGE no longer contains STANDARDS file segment
NEWS 13 Jul 22 USAN to be reloaded July 28, 2002;
saved answer sets no longer valid
NEWS 14 Jul 29 Enhanced polymer searching in REGISTRY
NEWS 15 Jul 30 NETFIRST to be removed from STN
NEWS 16 Aug 08 CANCERLIT reload
NEWS 17 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 18 Aug 08 NTIS has been reloaded and enhanced
NEWS 19 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE)
now available on STN
NEWS 20 Aug 19 IFIPAT, IFICDB, and IFIUDB have been reloaded
NEWS 21 Aug 19 The MEDLINE file segment of TOXCENTER has been reloaded
NEWS 22 Aug 26 Sequence searching in REGISTRY enhanced
NEWS 23 Sep 03 JAPIO has been reloaded and enhanced
NEWS 24 Sep 16 Experimental properties added to the REGISTRY file
NEWS 25 Sep 16 Indexing added to some pre-1967 records in CA/CAPLUS
NEWS 26 Sep 16 CA Section Thesaurus available in CAPLUS and CA
NEWS 27 Oct 01 CASREACT Enriched with Reactions from 1907 to 1985

NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d,
CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002
NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 17:57:47 ON 04 OCT 2002

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 17:57:55 ON 04 OCT 2002

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STRUCTURE FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5

DICTIONARY FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:

<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> list saved

LIST IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.

For a list of commands available to you in the current file, enter

"HELP COMMANDS" at an arrow prompt (=>).

=> saved

SAVED IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.

For a list of commands available to you in the current file, enter

"HELP COMMANDS" at an arrow prompt (=>).

=> help commands

Enter one of these commands at the arrow prompt (=>).

ACTIVATE ----- Assign L#s to saved query or answer set.
 ANALYZE ----- Build expand terms from answer fields.
 ARCHIVE ----- Purchase rights for archiving.
 BATCH ----- Request a batch search.
 DELETE ----- Delete saved or current session items.
 DISPLAY ----- Display saved or current session items.
 DOWNLOAD ----- Transfer data using Kermit protocol.
 EDIT ----- Modify the text of an E-number entry.
 EXPAND ----- Look at the index around a term.
 FILE ----- Specify the search and display file.
 HELP ----- For help on how to use the system.
 INDEX ----- Specify the Index environment.
 LOGOFF ----- End the online session.
 NEWS ----- Display current news about the system.
 ORDER ----- Order an original document or copy.
 POLYLINK ----- Search Using Polymer Linking
 PRINT ----- Print answers offline.
 QUERY ----- Define a search question (query).
 REDISTRIBUTE - Purchase rights for 'redistribution'.
 SAVE ----- Save an L-numbered query or answer set.
 SCREEN ----- Define a set of screens for searching.
 SDI ----- Request searches be run on file updates.
 SEARCH ----- Perform a search.
 SELECT ----- Build expand terms from answer fields.
 SEQLINK ----- Search Using Sequence Linking
 SET ----- Set terminal and interaction options.
 SORT ----- Sort Answers of an L-number list
 STRUCTURE ----- Create a structure for searching.
 TABULATE ----- Analyze SmartSELECT L-number data.
 TRANSFER ----- Extract terms from answers and search.
 ? ----- The same as HELP.

Enter one of these commands at the arrow prompt (=>).

For more detailed information about a command, enter HELP followed by the command name at an arrow prompt.

Example: => HELP SEARCH

For information on entering multiple commands at a single arrow prompt, enter HELP COMMAND STACKING.

=> display saved

NAME	CREATED	NOTES/TITLE
F403066/A	07 JUL 2000	3466 ANSWERS IN FILE REGISTRY
F485188/A	22 AUG 2000	15 ANSWERS IN FILE REGISTRY
F582302/A	29 MAR 2001	49 ANSWERS IN FILE REGISTRY
F9367871/A	15 MAY 2002	1856 ANSWERS IN FILE REGISTRY
F9386258/A	19 SEP 2000	179 ANSWERS IN FILE REGISTRY

F9396156/A	05 JUL 2000	28 ANSWERS IN FILE REGISTRY
F9445504/A	02 JUN 2000	63 ANSWERS IN FILE REGISTRY
F9446147/A	14 AUG 2000	301 ANSWERS IN FILE REGISTRY
F9446423/A	27 JUN 2000	107 ANSWERS IN FILE REGISTRY
F9462797/A	16 JUL 2000	13 ANSWERS IN FILE REGISTRY
F9462812/A	19 SEP 2000	0 ANSWERS IN FILE REGISTRY
F9482810/A	15 AUG 2000	226 ANSWERS IN FILE REGISTRY
F9486327/A	15 JUL 2000	51 ANSWERS IN FILE REGISTRY
F9486646/A	08 MAR 2001	12669 ANSWERS IN FILE REGISTRY
F9489271/A	02 JAN 2001	155 ANSWERS IN FILE REGISTRY
F9499699/A	01 JUN 2000	7 ANSWERS IN FILE REGISTRY
F9501758/A	11 OCT 2000	42 ANSWERS IN FILE REGISTRY
F9504183/A	26 JUN 2000	3 ANSWERS IN FILE REGISTRY
F9508240/A	16 DEC 2000	363 ANSWERS IN FILE REGISTRY
F9508796/A	12 JUL 2000	91 ANSWERS IN FILE REGISTRY
F9508989/A	27 FEB 2001	176 ANSWERS IN FILE REGISTRY
F9509831/A	07 SEP 2000	165 ANSWERS IN FILE REGISTRY
F9513147/A	15 JUN 2000	360 ANSWERS IN FILE REGISTRY
F9538800/A	29 DEC 2000	114 ANSWERS IN FILE REGISTRY
F954237/A	26 SEP 2000	7030 ANSWERS IN FILE REGISTRY
F9546555/A	21 JUL 2000	2536 ANSWERS IN FILE REGISTRY
F9552036/A	21 FEB 2001	116 ANSWERS IN FILE REGISTRY
F9552969/A	09 SEP 2000	4659 ANSWERS IN FILE REGISTRY
F9566002/A	25 JUL 2000	201 ANSWERS IN FILE REGISTRY
F9583924/A	23 DEC 2000	910 ANSWERS IN FILE REGISTRY
F9592336/A	23 AUG 2000	141 ANSWERS IN FILE REGISTRY
F9606921/A	25 OCT 2000	167 ANSWERS IN FILE REGISTRY
F9622086/A	10 JAN 2001	136 ANSWERS IN FILE REGISTRY
F9627766/A	28 NOV 2000	0 ANSWERS IN FILE REGISTRY
F9634801/A	14 OCT 2000	417 ANSWERS IN FILE REGISTRY
F9687241/A	08 JUN 2001	195 ANSWERS IN FILE REGISTRY
F9693896/A	31 JAN 2001	9183 ANSWERS IN FILE REGISTRY
F9696504/A	15 MAY 2001	4549 ANSWERS IN FILE REGISTRY
F9718223/A	25 MAY 2001	226 ANSWERS IN FILE REGISTRY
F9787672/L	14 JUN 2001	2 L-NUMBERS
F9792866/A	25 SEP 2001	1833 ANSWERS IN FILE REGISTRY
F9868992/A	25 JAN 2002	574 ANSWERS IN FILE REGISTRY
F9977944/A	17 JAN 2002	771 ANSWERS IN FILE REGISTRY
R9380742/A	23 JUN 2000	114 ANSWERS IN FILE REGISTRY
R9655115/A	01 MAR 2001	322 ANSWERS IN FILE REGISTRY

=> delete saved

DELETE ALL SAVED QUERIES, STRUCTURES, SCREEN SETS, ANSWER SETS AND L# LISTS?

(Y)/N:y

ALL SAVED SCREENS, STRUCTURES, QUERIES, ANSWER SETS, AND L# LISTS DELETED

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

1.14

1.35

FILE 'CAPLUS' ENTERED AT 17:59:49 ON 04 OCT 2002

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Hong Liu

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FILE COVERS 1907 - 4 Oct 2002 VOL 137 ISS 15
FILE LAST UPDATED: 3 Oct 2002 (20021003/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> s wo9740049/pn
L1 1 WO9740049/PN

=> select l1 rn 1-
E1 THROUGH E197 ASSIGNED

=> fil reg		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	2.11	3.46

FILE 'REGISTRY' ENTERED AT 18:00:27 ON 04 OCT 2002
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STRUCTURE FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5
DICTIONARY FILE UPDATES: 3 OCT 2002 HIGHEST RN 459123-02-5

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> s el-el97

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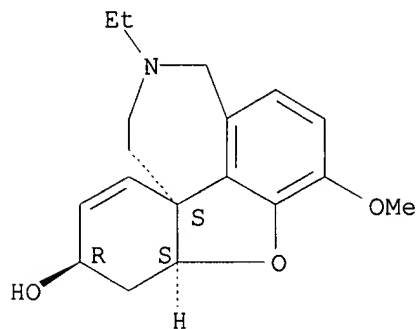
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1 35196-48-6/BI
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1 357-70-0/BI
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1 3647-69-6/BI
(3647-69-6/RN)
1 3878-55-5/BI
(3878-55-5/RN)
1 41303-52-0/BI
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1 4530-20-5/BI
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1 4584-46-7/BI
(4584-46-7/RN)
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1 60384-53-4/BI
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1 624-84-0/BI
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1 65232-62-4/BI
(65232-62-4/RN)
1 7250-67-1/BI
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1 7536-58-5/BI
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1 7803-57-8/BI
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1 79-07-2/BI
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OR
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110-53-2/BI OR 111-25-1/BI OR 111-36-4/BI OR 112-67-4/BI OR
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OR 179108-06-6/BI OR 179239-41-9/BI OR 180854-29-9/BI OR
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198987-86-9/BI OR 198987-87-0/BI OR 198987-88-1/BI OR
198987-89-
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198987-93-8/BI OR 198987-94-9/BI OR 198987-95-0/B

=> d scan

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-ethyl-4a,5,9,10,11,12-
hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H23 N O3

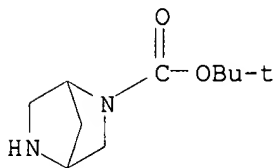
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

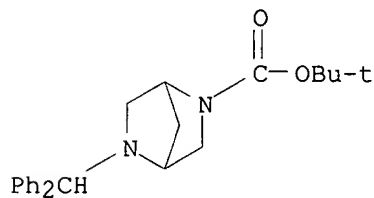
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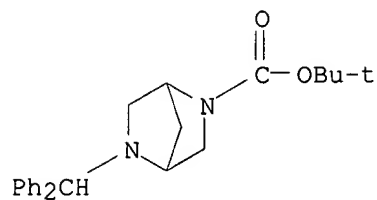
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 1,1-dimethylethyl ester (9CI)
 MF C10 H18 N2 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

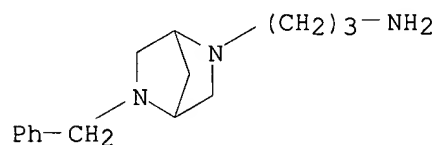
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 5-(diphenylmethyl)-, 1,1-dimethylethyl ester (9CI)
 MF C23 H28 N2 O2





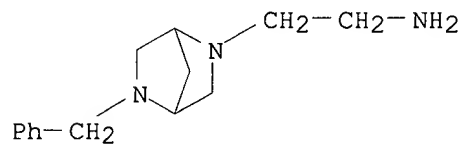
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-propanamine, 5-(phenylmethyl)- (9CI)
 MF C15 H23 N3



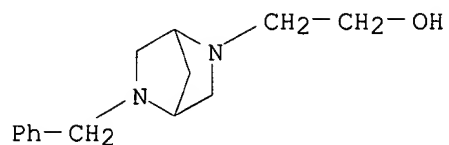
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-ethanamine, 5-(phenylmethyl)- (9CI)
 MF C14 H21 N3



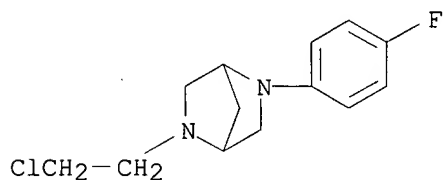
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-ethanol, 5-(phenylmethyl)- (9CI)
 MF C14 H20 N2 O



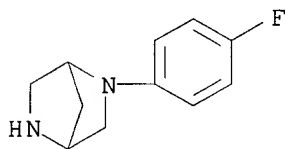
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-chloroethyl)-5-(4-fluorophenyl)- (9CI)
 MF C13 H16 Cl F N2



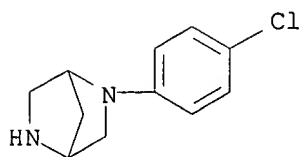
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-fluorophenyl)- (9CI)
 MF C11 H13 F N2



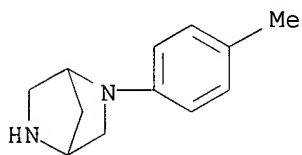
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)- (9CI)
 MF C11 H13 Cl N2



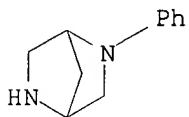
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-methylphenyl)- (9CI)
MF C12 H16 N2



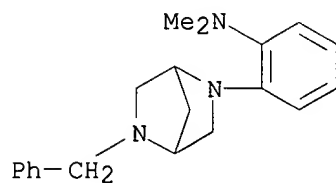
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-phenyl- (9CI)
MF C11 H14 N2
CI COM



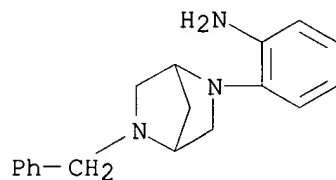
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenamine,
N,N-dimethyl-2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-
2-yl]- (9CI)
MF C20 H25 N3



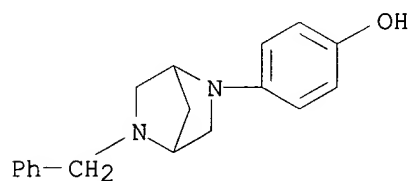
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenamine, 2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
 MF C18 H21 N3



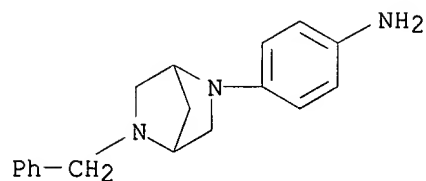
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Phenol, 4-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
 MF C18 H20 N2 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

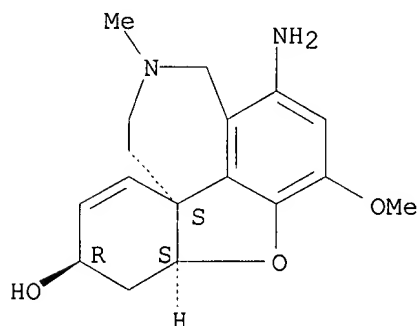
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenamine, 4-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]- (9CI)
 MF C18 H21 N3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-amino-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
 MF C17 H22 N2 O3

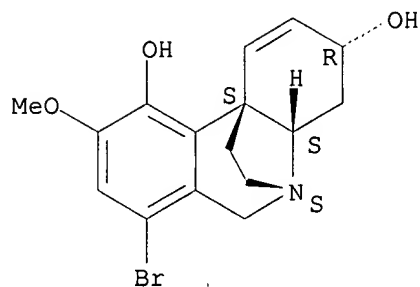
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 3H,6H-5,10b-Ethanophenanthridine-3,10-diol,
 7-bromo-4,4a-dihydro-9-methoxy-
 , (3.alpha.,4a.beta.,5.beta.,10b.beta.)- (9CI)
 MF C16 H18 Br N O3

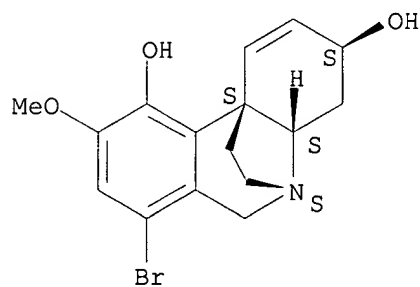
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 3H,6H-5,10b-Ethanophenanthridine-3,10-diol,
 7-bromo-4,4a-dihydro-9-methoxy-
 , (3.alpha.,4a.alpha.,5.alpha.,10b.alpha.)- (9CI)
 MF C16 H18 Br N O3

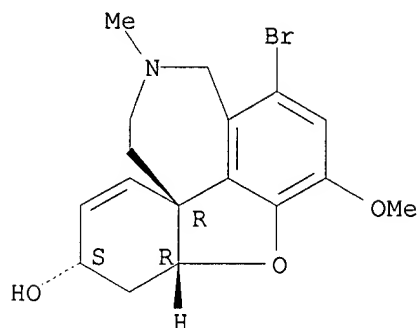
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-methyl-, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C17 H20 Br N O3

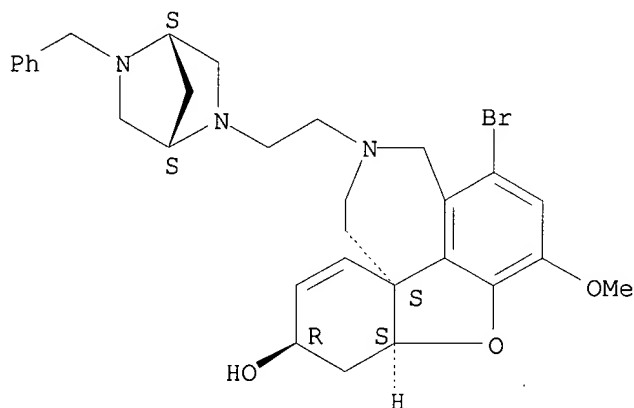
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-[2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-
 yl]ethyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
 MF C30 H36 Br N3 O3

Relative stereochemistry.

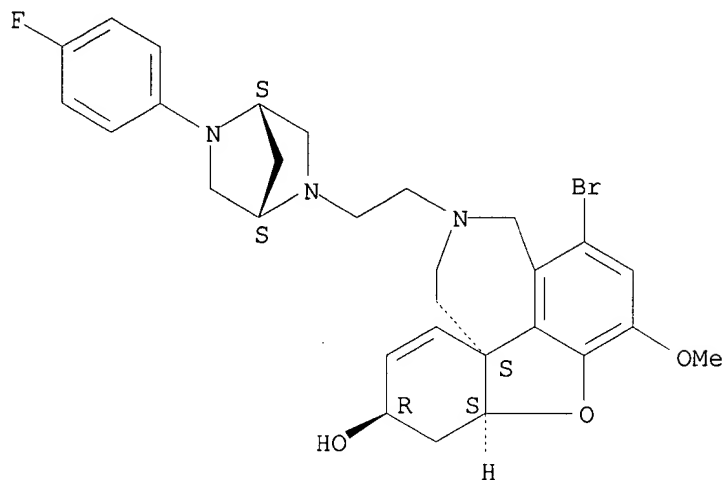


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-[2-[5-(4-
 fluorophenyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]ethyl]-4a,5,9,10,11,12-
 hexahydro-3-methoxy-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)

MF C29 H33 Br F N3 O3

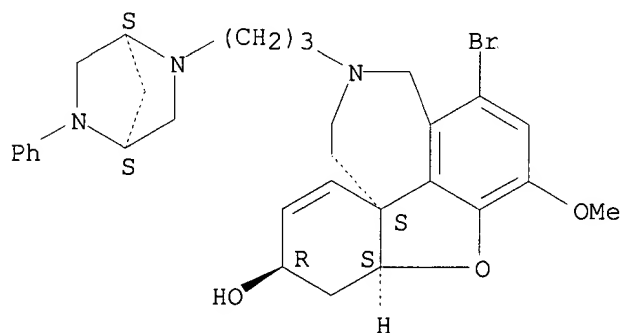
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-[3-(5-phenyl-2,5-diazabicyclo[2.2.1]hept-2-yl)propyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
 MF C30 H36 Br N3 O3

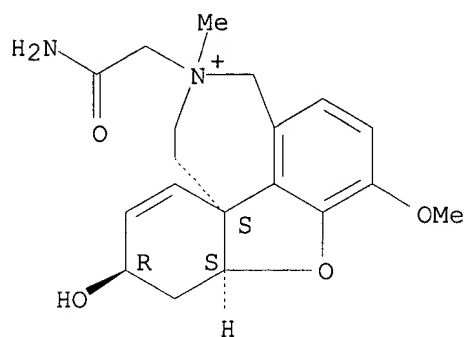
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 11-(2-amino-2-oxoethyl)-
4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-, bromide,
(4aS,6R,8aS)-[partial]- (9CI)
MF C19 H25 N2 O4 . Br

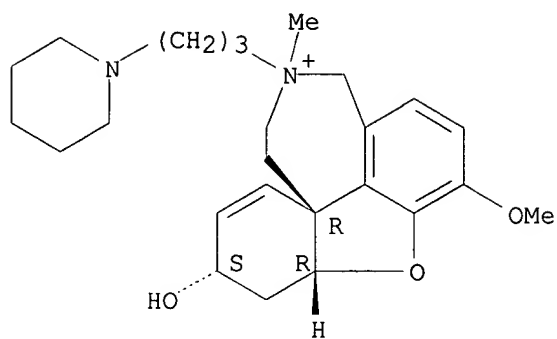
Absolute stereochemistry.



● Br⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-
hydroxy-3-methoxy-11-methyl-11-[3-(1-piperidinyl)propyl]-, chloride,
(4aR,6S,8aR)-[partial]- (9CI)
MF C25 H37 N2 O3 . Cl

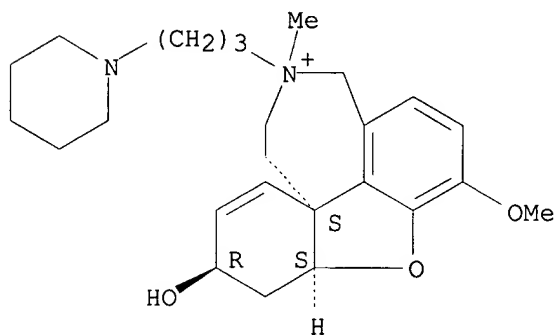
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[3-(1-piperidinyl)propyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
 MF C25 H37 N2 O3 . Cl

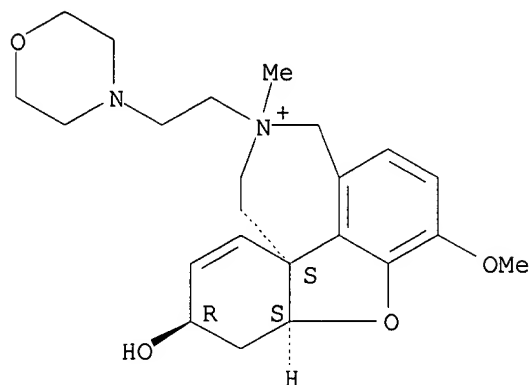
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(4-morpholinyl)ethyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
 MF C23 H33 N2 O4 . Cl

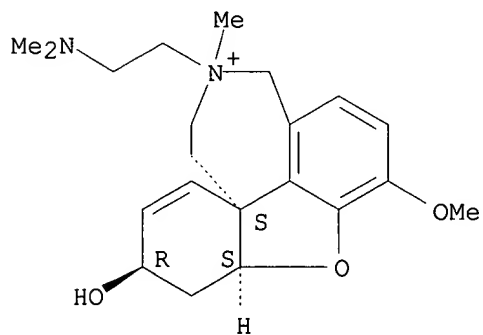
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 11-[2-(dimethylamino)ethyl]-
4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-, chloride,
(4aS,6R,8aS)-[partial]- (9CI)
MF C21 H31 N2 O3 . Cl

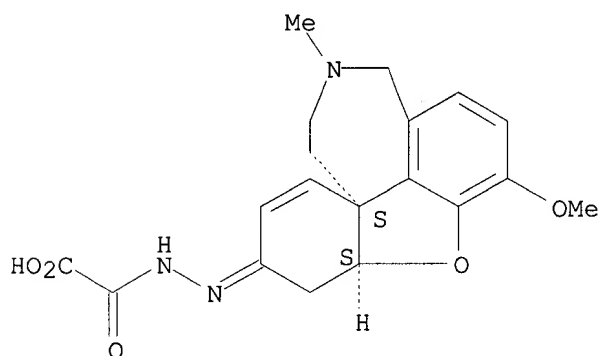
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanedioic acid, mono[(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)hydrazide], (4aR*,8aR*)-(9CI)
MF C19 H21 N3 O5

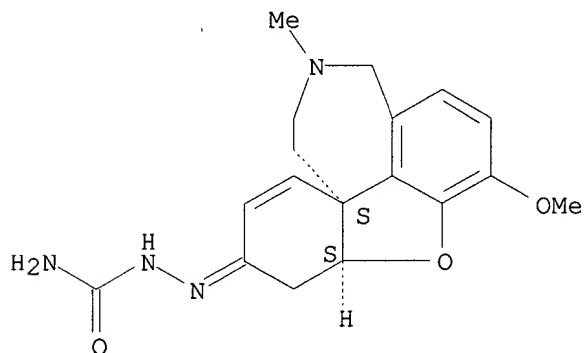
Relative stereochemistry.
Double bond geometry unknown.

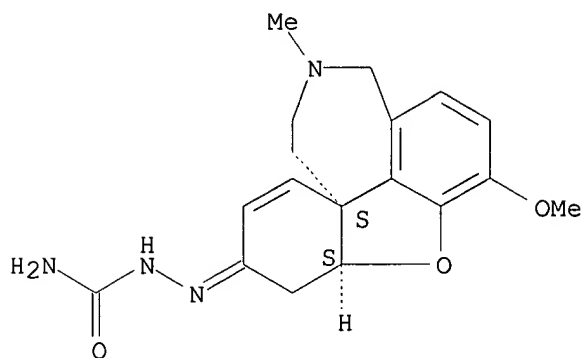


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxamide,
2-[(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-], (4aR*,8aR*)-(9CI)
MF C18 H22 N4 O3

Relative stereochemistry.
Double bond geometry unknown.

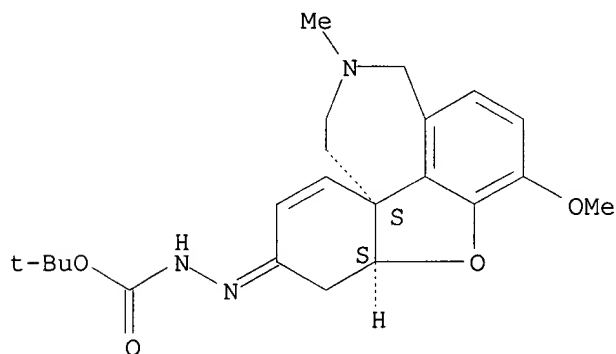




PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hydrazinecarboxylic acid, (4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, 1,1-dimethylethyl ester,
 (4aR*,8aR*)- (9CI)
 MF C22 H29 N3 O4

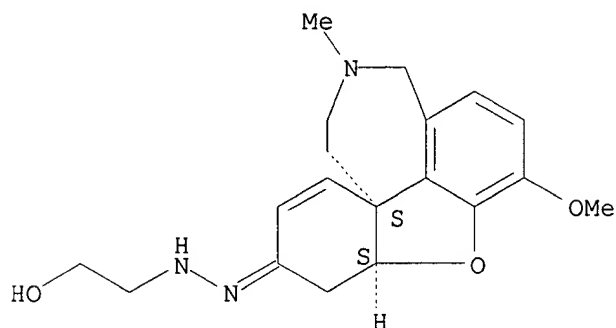
Relative stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (2-hydroxyethyl)hydrazone, [4aS-(4aR*,8aR*)]- (9CI)
 MF C19 H25 N3 O3

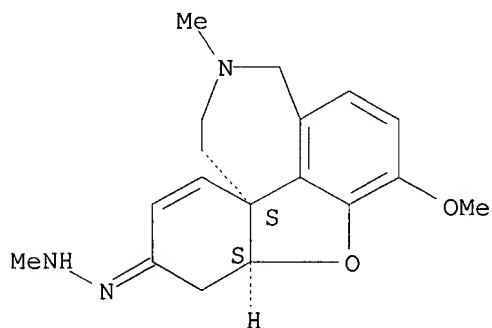
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
methoxy-11-methyl-, methylhydrazone, [4aS-(4aR*,8aR*)]- (9CI)
MF C18 H23 N3 O2

Absolute stereochemistry.
Double bond geometry unknown.

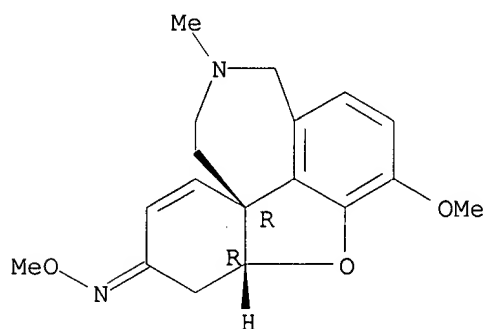


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-

methoxy-11-methyl-, O-methyloxime, [4aR-(4aR*,8aR*)]- (9CI)
MF C18 H22 N2 O3

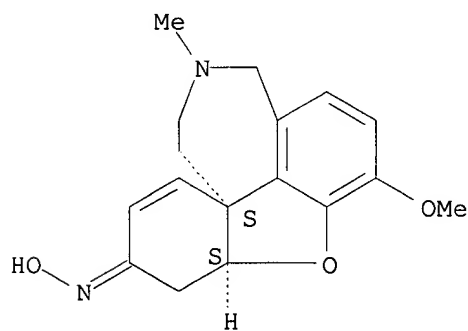
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, oxime, [4aS-(4aR*,8aR*)]- (9CI)
MF C17 H20 N2 O3

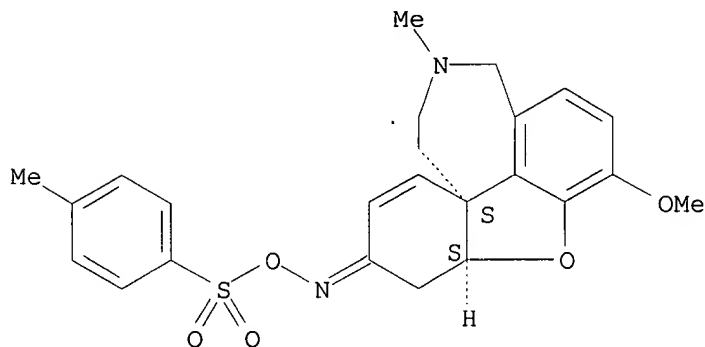
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, O-[(4-methylphenyl)sulfonyl]oxime, (4aR*,8aR*)- (9CI)
MF C24 H26 N2 O5 S

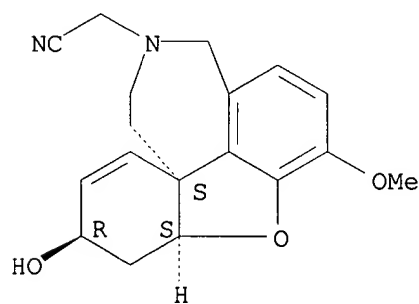
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetonitrile, 4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C18 H20 N2 O3

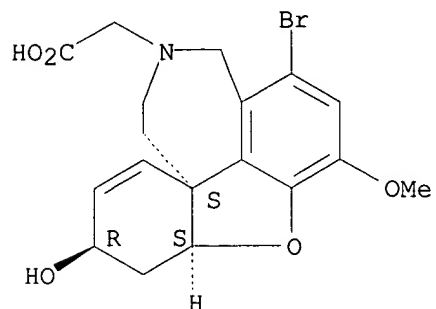
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O5

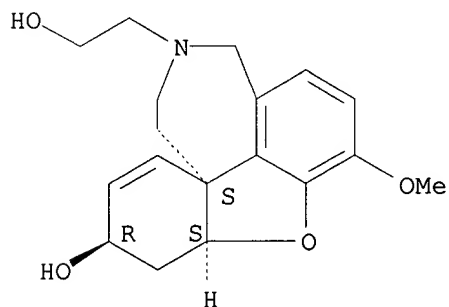
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-ethanol,
4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C18 H23 N O4

Relative stereochemistry.

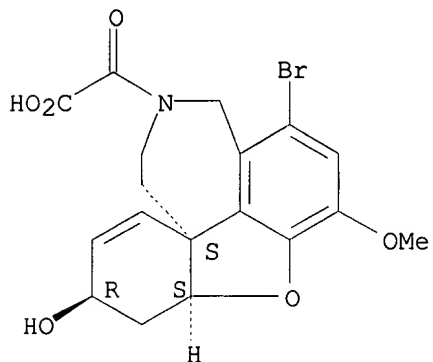


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.alpha.-oxo-,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H18 Br N O6

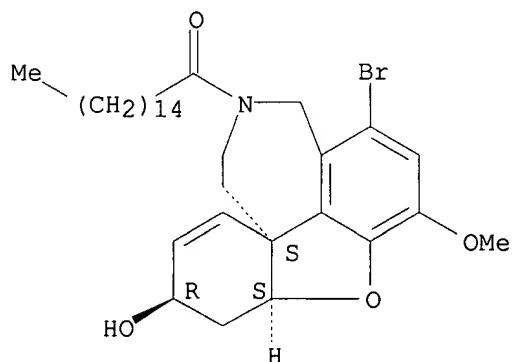
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
hexahydro-3-methoxy-11-(1-oxohexadecyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C32 H48 Br N O4

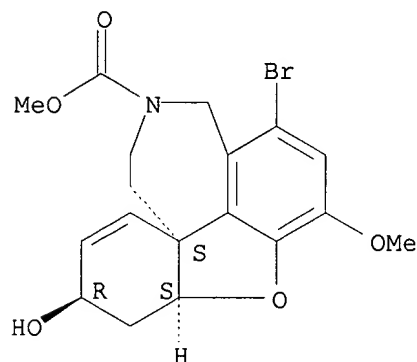
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxylic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, methyl ester,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O5

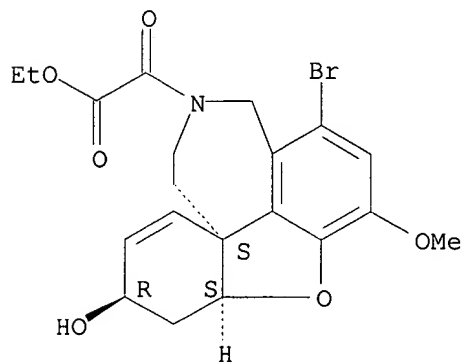
Relative stereochemistry.



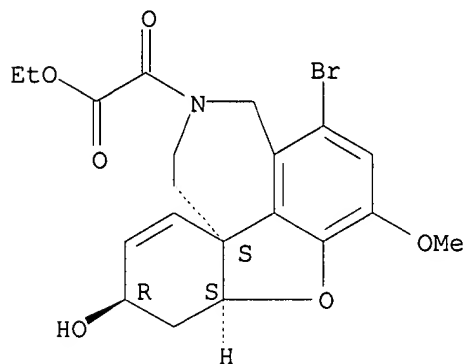
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.alpha.-oxo-, ethyl
ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H22 Br N O6

Relative stereochemistry.



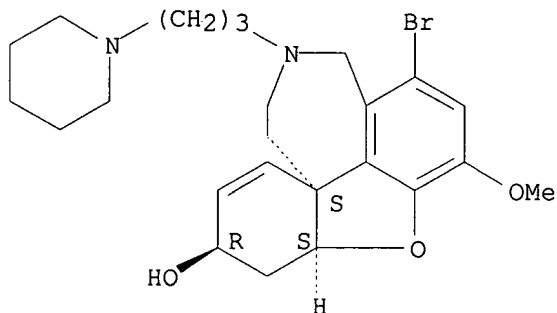
Hong Liu



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-[3-(1-piperidinyl)propyl]-,
 (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C24 H33 Br N2 O3

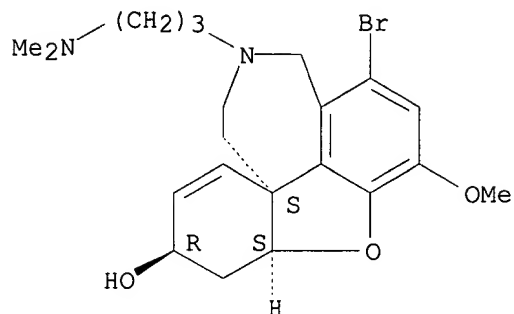
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-[3-
 (dimethylamino)propyl]-4a,5,9,10,11,12-hexahydro-3-methoxy-,
 (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C21 H29 Br N2 O3

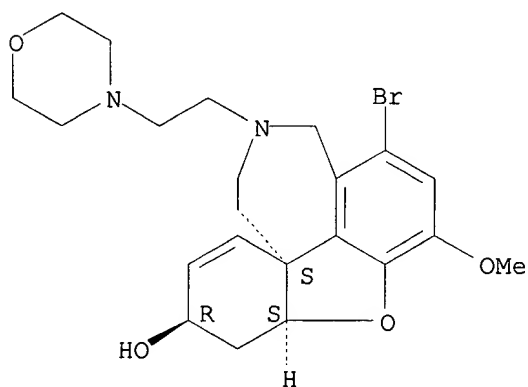
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-[2-(4-morpholinyl)ethyl]-,
 (4a.alpha.,6.beta.,8aR*)-
 (9CI)
 MF C22 H29 Br N2 O4

Relative stereochemistry.

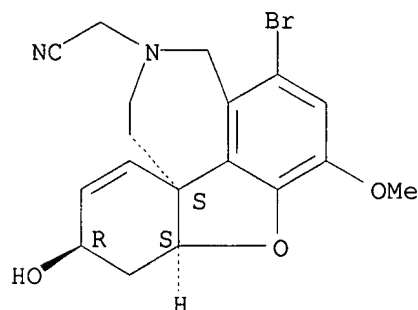


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetonitrile,
 1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-,
 (4a.alpha.,6.beta.,8aR*)- (9CI)

MF C18 H19 Br N2 O3

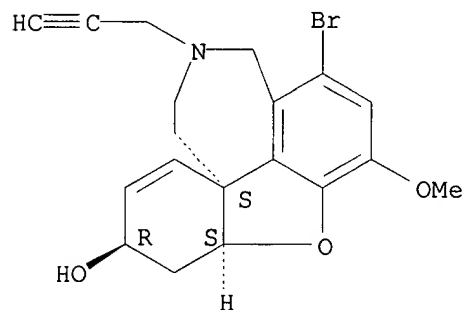
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-(2-propynyl)-, (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C19 H20 Br N O3

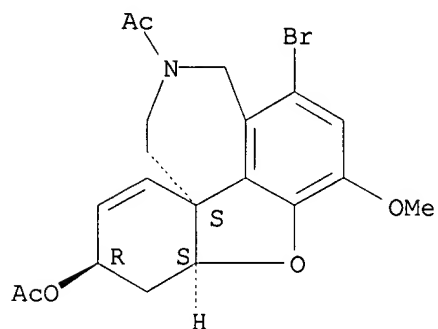
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-acetyl-1-bromo-
 4a,5,9,10,11,12-hexahydro-3-methoxy-, acetate (ester),
 (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C20 H22 Br N O5

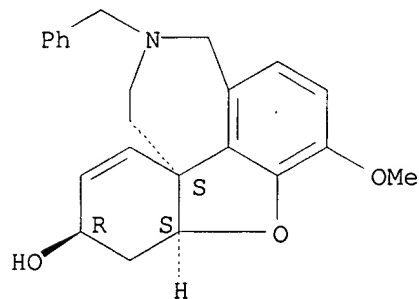
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C23 H25 N O3

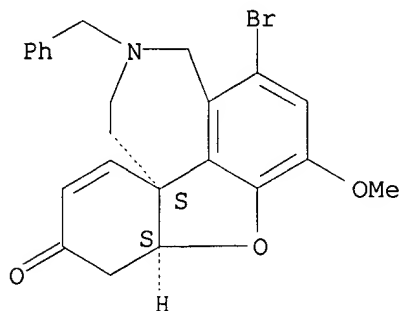
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-, (4aR*,8aR*)- (9CI)
 MF C23 H22 Br N O3

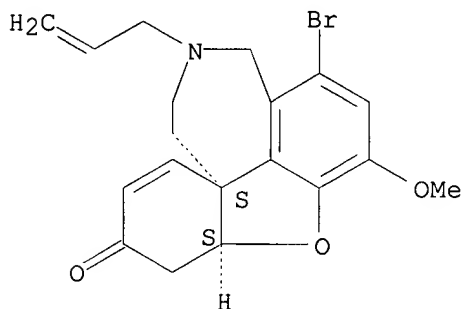
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-(2-propenyl)-, (4aR*,8aR*)- (9CI)
 MF C19 H20 Br N O3

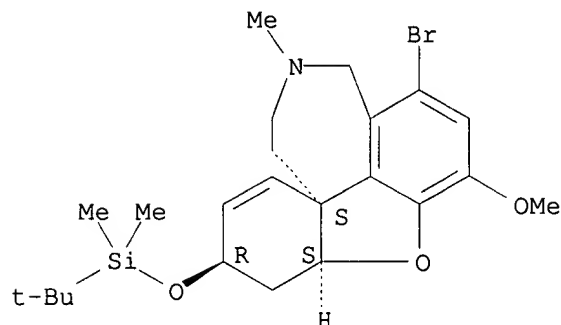
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 1-bromo-6-[[[(1,1-
 dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-11-
 methyl-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C23 H34 Br N O3 Si

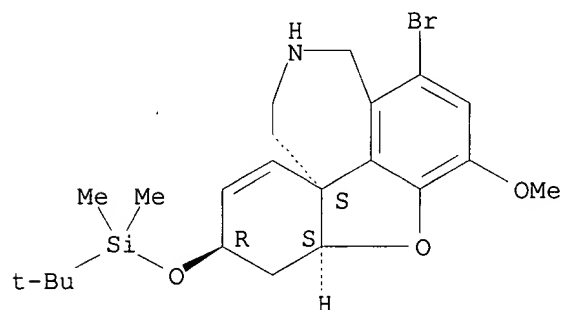
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 1-bromo-6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C22 H32 Br N O3 Si

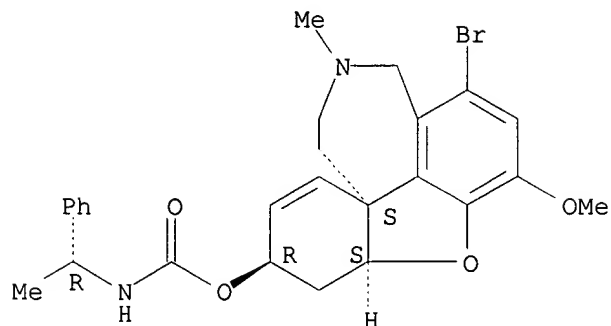
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Carbamic acid, (1-phenylethyl)-, 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4a.alpha.,6.beta.(S*),8aR*]- (9CI)
 MF C26 H29 Br N2 O4

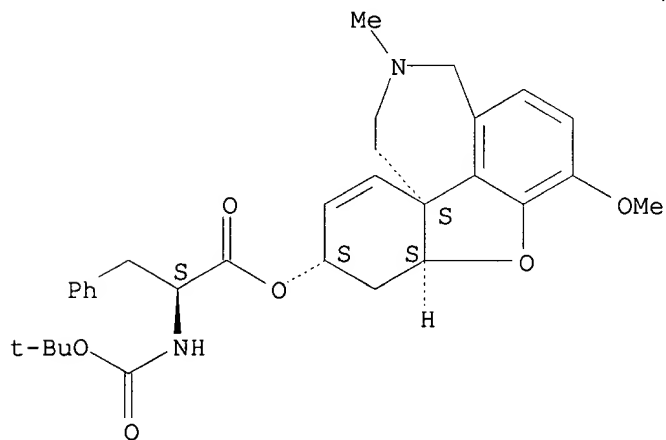
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN L-Phenylalanine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-
 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
 ef][2]benzazepin-6-yl ester (9CI)
 MF C31 H38 N2 O6

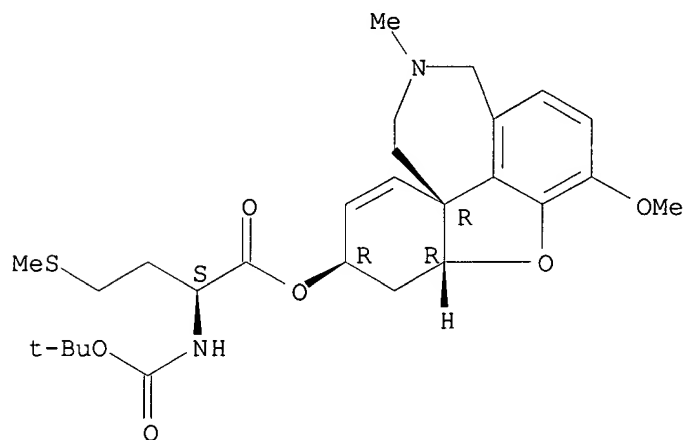
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aR,6R,8aR)-
4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
ef][2]benzazepin-6-yl ester (9CI)
MF C27 H38 N2 O6 S

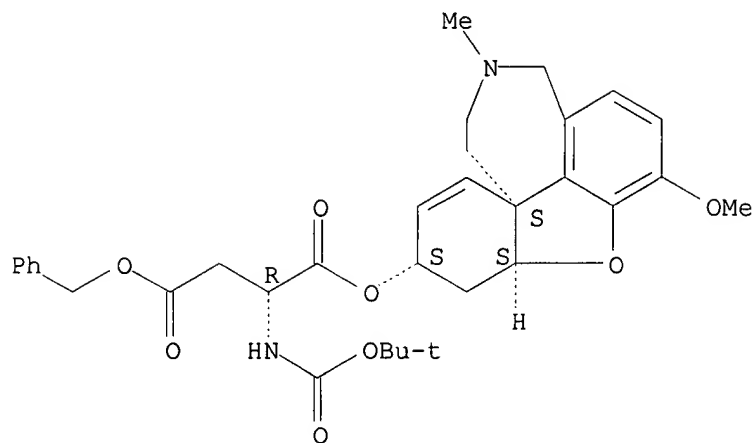
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN D-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aS,6S,8aS)-
4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

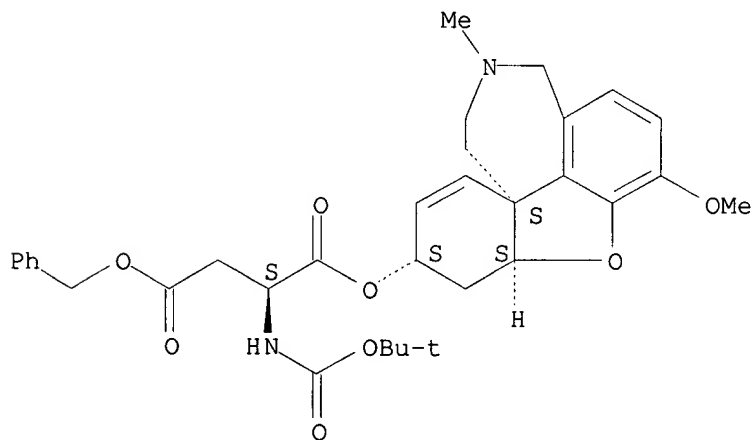
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aS,6S,8aS)-
4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

Absolute stereochemistry.

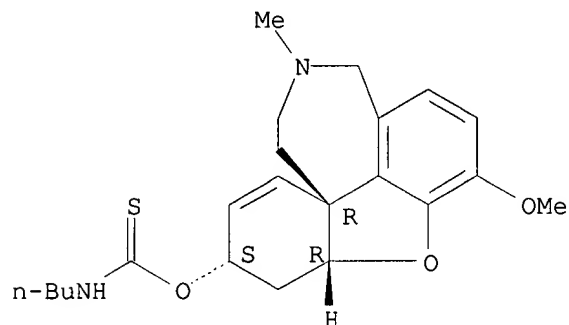


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

Hong Liu

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Carbamothioic acid, butyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester,
 [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C22 H30 N2 O3 S

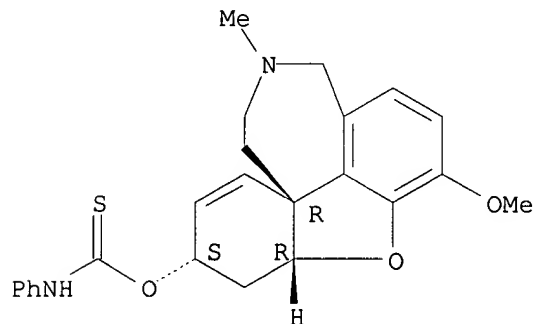
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Carbamothioic acid, phenyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester,
 [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C24 H26 N2 O3 S

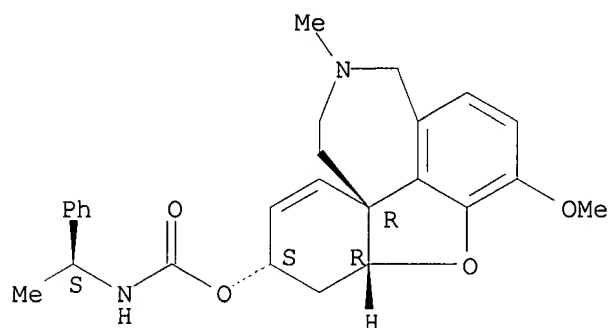
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester,
[4aR-[4a.alpha.,6.beta.(S*),8aR*]]- (9CI)
MF C26 H30 N2 O4

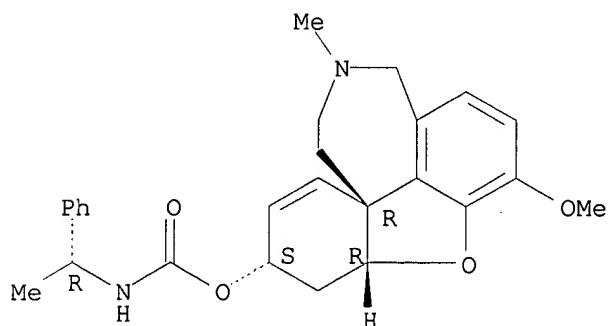
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester,
[4aR-[4a.alpha.,6.beta.(R*),8aR*]]- (9CI)
MF C26 H30 N2 O4

Absolute stereochemistry.

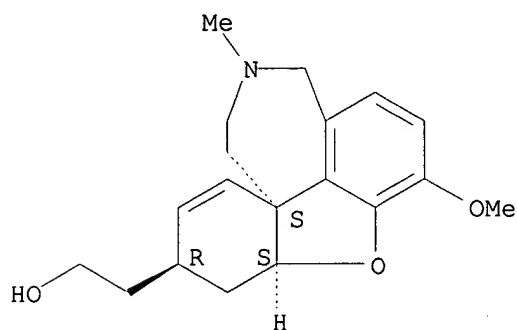


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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-6-ethanol,
4a,5,9,10,11,12-hexahydro-
3-methoxy-11-methyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C19 H25 N O3

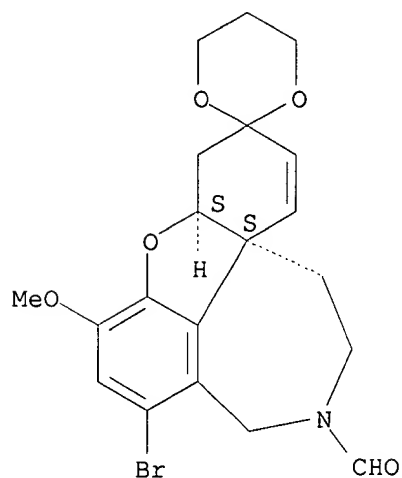
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxane]-11(12H)-
carboxaldehyde, 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-, (4aR*,8aR*)-
(9CI)
MF C20 H22 Br N O5

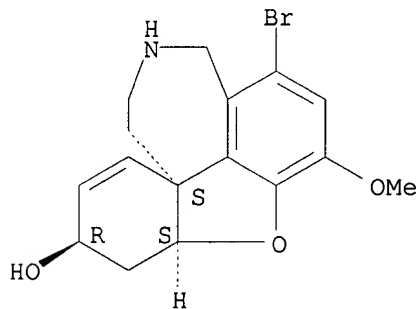
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-, (4aS,6R,8aS)- (9CI)
 MF C16 H18 Br N O3

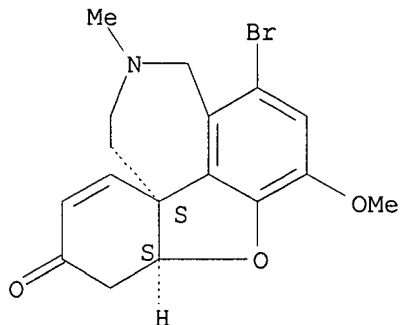
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
 MF C17 H18 Br N O3

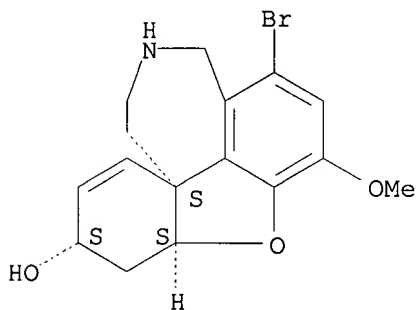
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-, (4aR,6R,8aR)-rel- (9CI)
 MF C16 H18 Br N O3

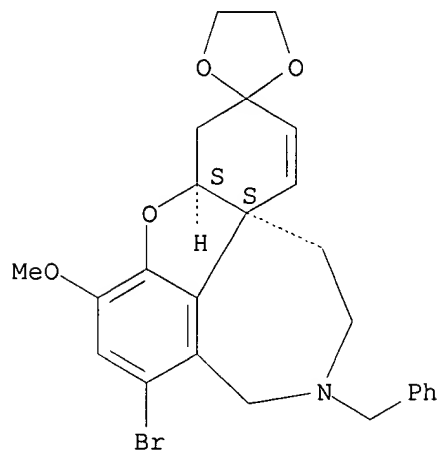
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane],
 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-(phenylmethyl)-,
 (4aR*,8aR*)- (9CI)
 MF C25 H26 Br N O4

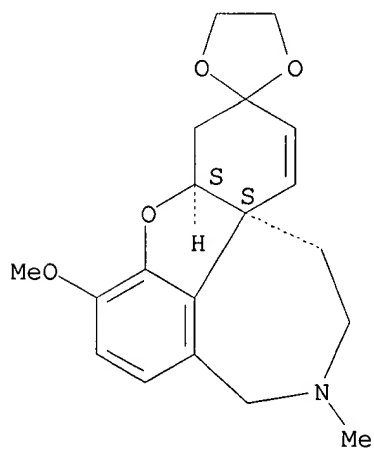
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane],
 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
 MF C19 H23 N O4

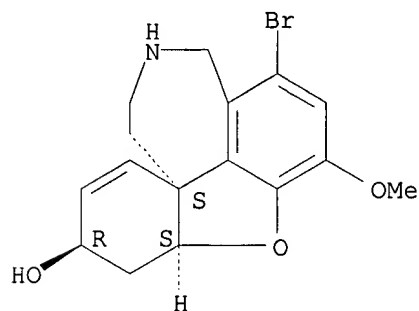
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
hexahydro-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
MF C16 H18 Br N O3

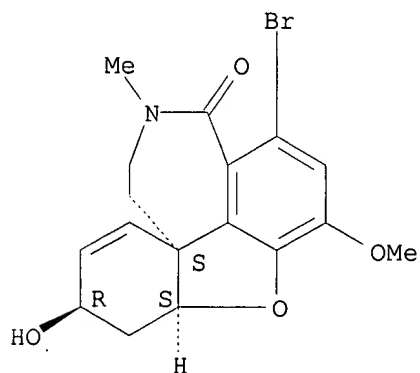
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

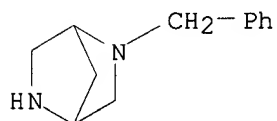
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-12(9H)-one, 1-bromo-4a,5,10,11-
tetrahydro-6-hydroxy-3-methoxy-11-methyl-, (4aR,6S,8aR)-rel- (9CI)
MF C17 H18 Br N O4

Relative stereochemistry.



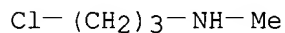
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(phenylmethyl)-, dihydrobromide (9CI)
MF C12 H16 N2 . 2 Br H



● 2 HBr

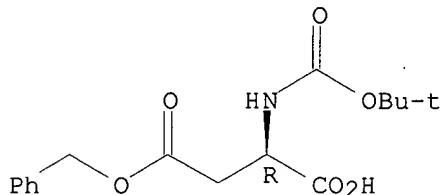
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Propanamine, 3-chloro-N-methyl- (9CI)
MF C4 H10 Cl N
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

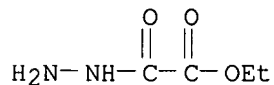
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN D-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 4-(phenylmethyl)
ester (9CI)
MF C16 H21 N O6

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

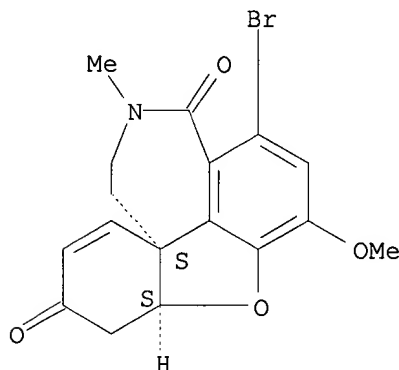
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanedioic acid, monoethyl ester, hydrazide (9CI)
MF C4 H8 N2 O3
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Galanthamine, 8-bromo-3-deoxy-3,9-dioxo- (9CI)
MF C17 H16 Br N O4

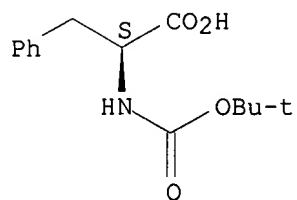
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Phenylalanine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C14 H19 N O4
CI COM

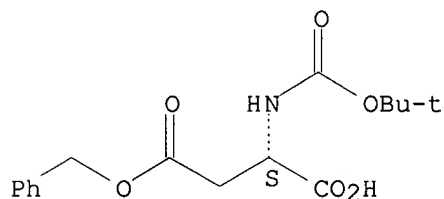
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

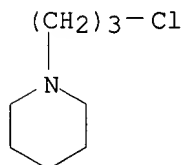
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 4-(phenylmethyl)
 ester
 (9CI)
 MF C16 H21 N O6
 CI COM

Absolute stereochemistry.



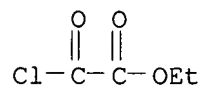
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Piperidine, 1-(3-chloropropyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
 MF C8 H16 Cl N . Cl H



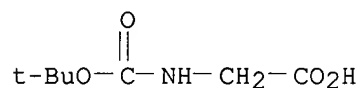
● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, chlorooxo-, ethyl ester (9CI)
MF C4 H5 Cl O3



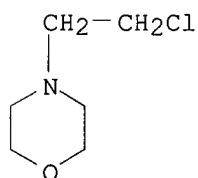
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Glycine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C7 H13 N O4
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

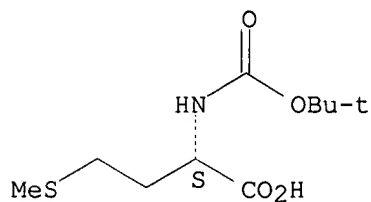
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Morpholine, 4-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C6 H12 Cl N O . Cl H



● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C10 H19 N O4 S
CI COM

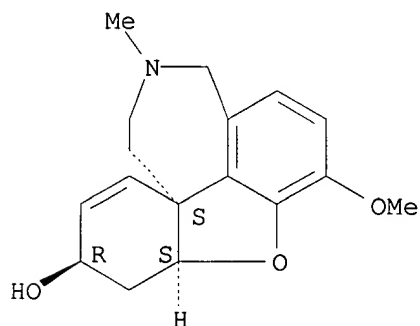
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

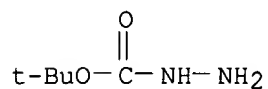
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, hydrobromide, (4aS,6R,8aS)- (9CI)
 MF C17 H21 N O3 . Br H
 CI COM

Absolute stereochemistry. Rotation (-).



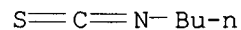
● HBr

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hydrazinecarboxylic acid, 1,1-dimethylethyl ester (9CI)
 MF C5 H12 N2 O2
 CI COM



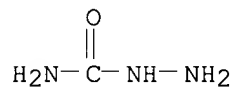
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Butane, 1-isothiocyanato- (9CI)
 MF C5 H9 N S
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

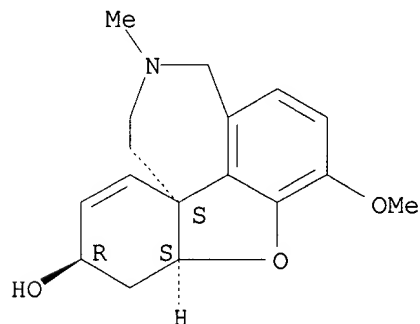
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hydrazinecarboxamide, monohydrochloride (9CI)
 MF C H5 N3 O . Cl H
 CI COM



● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
 MF C17 H21 N O3
 CI COM

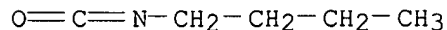
Absolute stereochemistry. Rotation (-).



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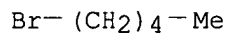
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Butane, 1-isocyanato- (9CI)
MF C5 H9 N O
CI COM



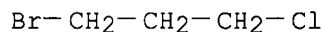
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Pentane, 1-bromo- (6CI, 8CI, 9CI)
MF C5 H11 Br
CI COM



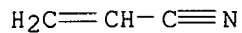
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Propane, 1-bromo-3-chloro- (6CI, 7CI, 8CI, 9CI)
MF C3 H6 Br Cl
CI COM



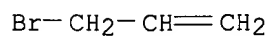
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2-Propenenitrile (9CI)
MF C3 H3 N
CI COM



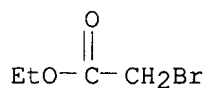
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1-Propene, 3-bromo- (9CI)
MF C3 H5 Br
CI COM



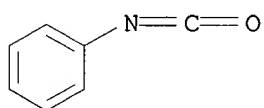
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, bromo-, ethyl ester (6CI, 8CI, 9CI)
MF C4 H7 Br O2
CI COM



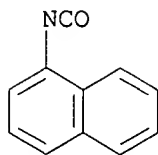
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, isocyanato- (9CI)
MF C7 H5 N O
CI COM



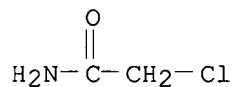
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Naphthalene, 1-isocyanato- (9CI)
MF C11 H7 N O
CI COM



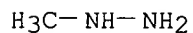
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Acetamide, 2-chloro- (6CI, 7CI, 8CI, 9CI)
 MF C2 H4 Cl N O
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

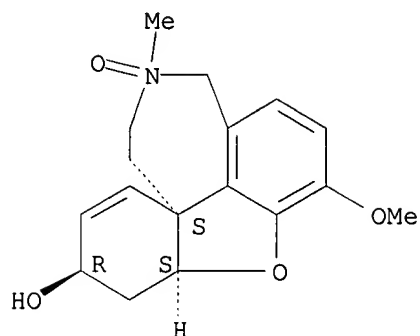
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hydrazine, methyl- (6CI, 8CI, 9CI)
 MF C H6 N2
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

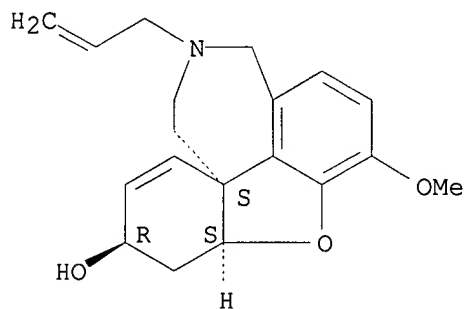
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, 11-oxide, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C17 H21 N O4

Absolute stereochemistry.



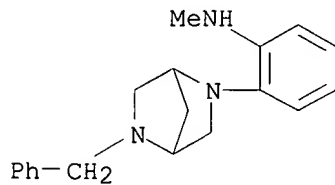
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(2-propenyl)-, (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C19 H23 N O3

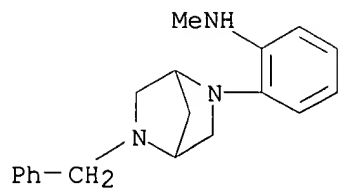
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

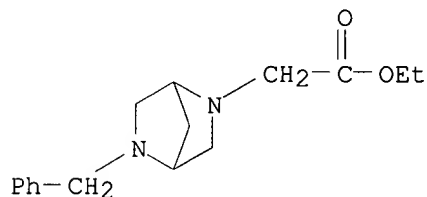
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenamine, N-methyl-2-[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-yl]-(9CI)
 MF C19 H23 N3





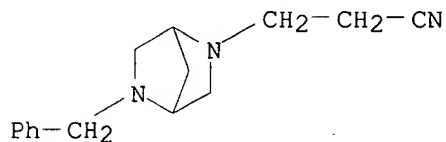
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-acetic acid, 5-(phenylmethyl)-, ethyl
 ester (9CI)
 MF C16 H22 N2 O2



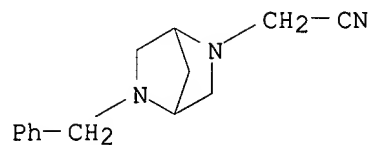
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-propanenitrile, 5-(phenylmethyl)- (9CI)
 MF C15 H19 N3



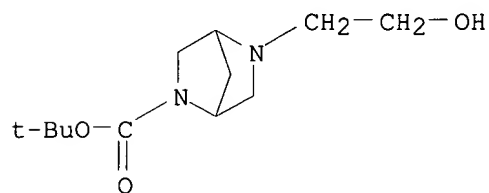
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-acetonitrile, 5-(phenylmethyl)- (9CI)
 MF C14 H17 N3



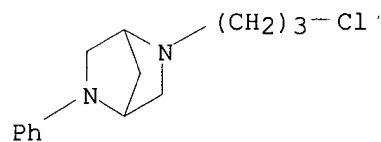
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane-2-carboxylic acid, 5-(2-hydroxyethyl)-,
 1,1-dimethylethyl ester (9CI)
 MF C12 H22 N2 O3



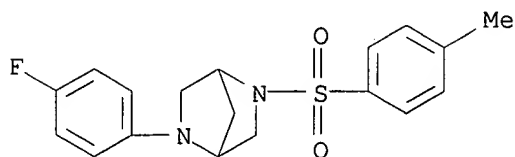
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(3-chloropropyl)-5-phenyl- (9CI)
 MF C14 H19 Cl N2



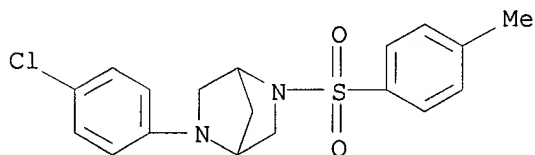
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-fluorophenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
 MF C18 H19 F N2 O2 S



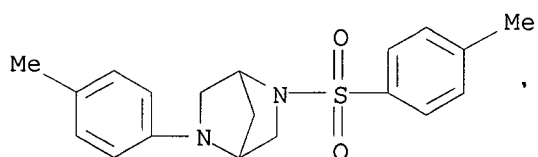
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
 MF C18 H19 Cl N2 O2 S



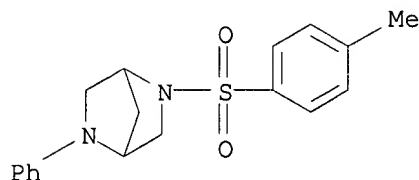
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-methylphenyl)-5-[(4-methylphenyl)sulfonyl]- (9CI)
 MF C19 H22 N2 O2 S



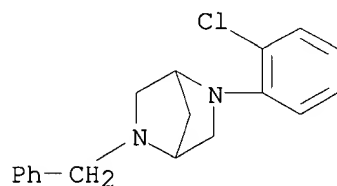
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-[(4-methylphenyl)sulfonyl]-5-phenyl- (9CI)
 MF C18 H20 N2 O2 S



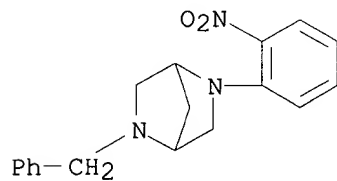
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-chlorophenyl)-5-(phenylmethyl)-
(9CI)
MF C18 H19 Cl N2



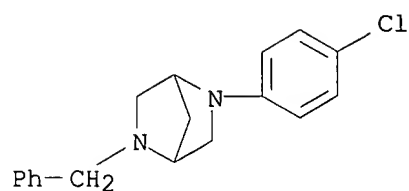
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(2-nitrophenyl)-5-(phenylmethyl)- (9CI)
MF C18 H19 N3 O2



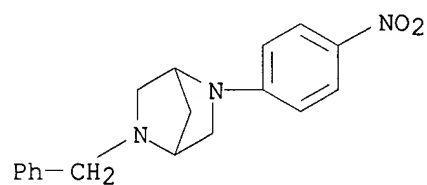
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-chlorophenyl)-5-(phenylmethyl)-
(9CI)
MF C18 H19 Cl N2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

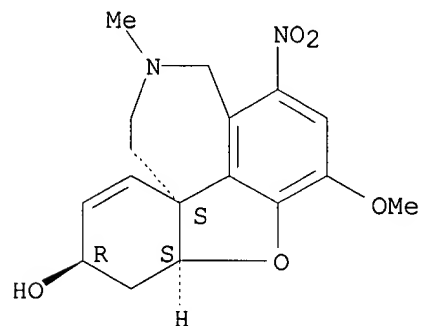
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2,5-Diazabicyclo[2.2.1]heptane, 2-(4-nitrophenyl)-5-(phenylmethyl)- (9CI)
 MF C18 H19 N3 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-1-nitro-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C17 H20 N2 O5

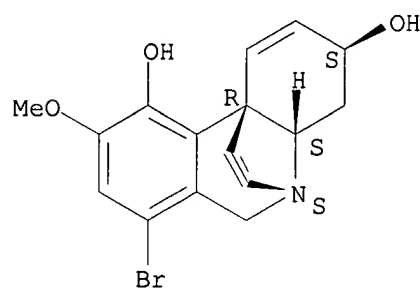
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

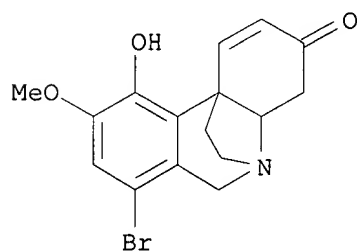
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 3H,6H-5,10b-Ethynophenanthridine-3,10-diol,
 7-bromo-4,4a-dihydro-9-methoxy-
 , (3.alpha.,4a.alpha.,5.alpha.,10b.alpha.)- (9CI)
 MF C16 H14 Br N O3

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 3H,6H-5,10b-Ethanophenanthridin-3-one, 7-bromo-4,4a-dihydro-10-hydroxy-9-
 methoxy- (9CI)
 MF C16 H16 Br N O3

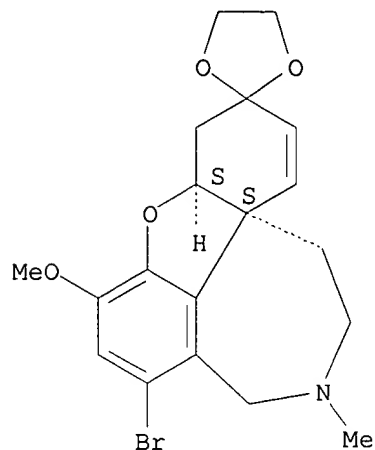


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane],
 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)-
 (9CI)

MF C19 H22 Br N O4

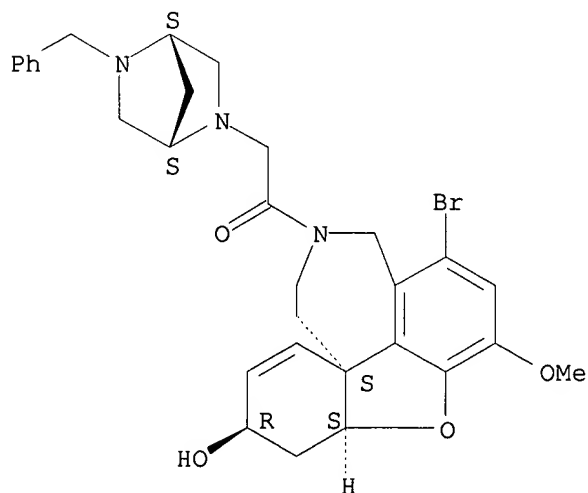
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
hexahydro-3-methoxy-11-[[5-(phenylmethyl)-2,5-diazabicyclo[2.2.1]hept-2-
yl]acetyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H34 Br N3 O4

Relative stereochemistry.

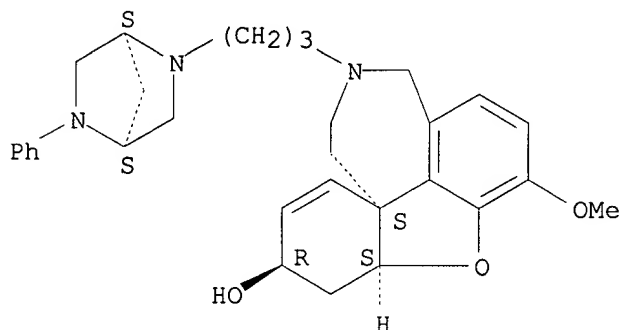


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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-[3-(5-phenyl-2,5-diazabicyclo[2.2.1]hept-2-yl)propyl]-, [4a.alpha.,6.beta.,8aR*,11(1R*,4R*)]- (9CI)
MF C30 H37 N3 O3

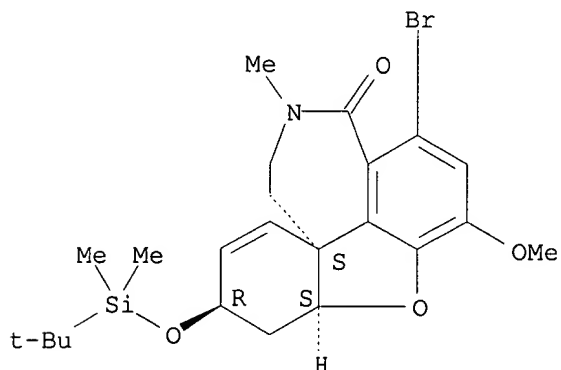
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-12(9H)-one, 1-bromo-6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,10,11-tetrahydro-3-methoxy-11-methyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C23 H32 Br N O4 Si

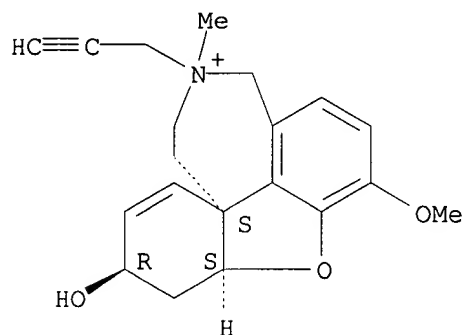
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

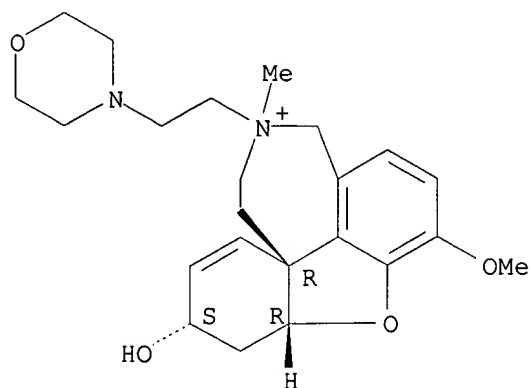
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-(2-propynyl)-, bromide, (4aS,6R,8aS)-[partial]- (9CI)
 MF C20 H24 N O3 . Br

Absolute stereochemistry.



L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(4-morpholinyl)ethyl]-, chloride, (4aR,6S,8aR)-[partial]- (9CI)
 MF C23 H33 N2 O4 . Cl

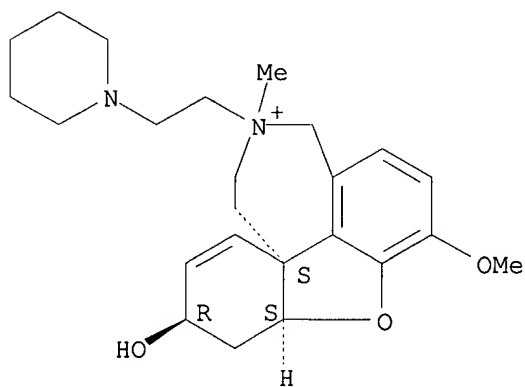
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-hydroxy-3-methoxy-11-methyl-11-[2-(1-piperidinyl)ethyl]-, chloride, (4aS,6R,8aS)-[partial]- (9CI)
 MF C24 H35 N2 O3 . Cl

Absolute stereochemistry.

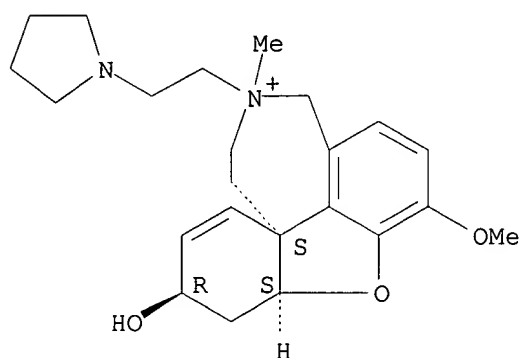


● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepinium, 4a,5,9,10,11,12-hexahydro-6-

hydroxy-3-methoxy-11-methyl-11-[2-(1-pyrrolidinyl)ethyl]-, chloride,
 (4a*S*,6*R*,8a*S*)-[partial]- (9*CI*)
 MF C23 H33 N2 O3 . Cl

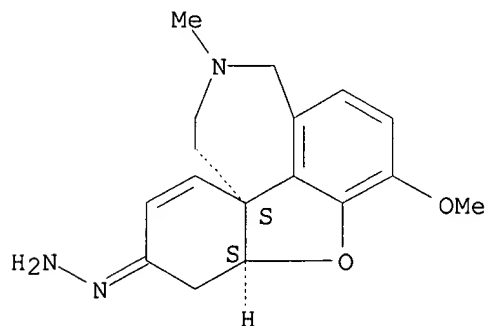
Absolute stereochemistry.



● Cl⁻

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
 methoxy-11-methyl-, hydrazone, [4a*S*-(4a*R**,8a*R**)]- (9*CI*)
 MF C17 H21 N3 O2

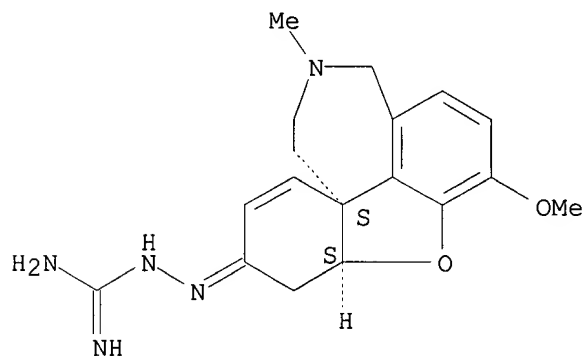
Relative stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboximidamide,
2-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-
6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, (4aR*,8aR*)- (9CI)
MF C18 H23 N5 O2

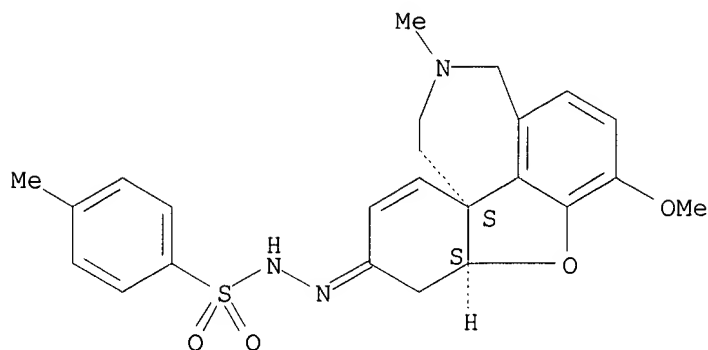
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzenesulfonic acid, 4-methyl-, (4a,5,9,10,11,12-hexahydro-3-methoxy-11-
methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)hydrazide,
(4aR*,8aR*)- (9CI)
MF C24 H27 N3 O4 S

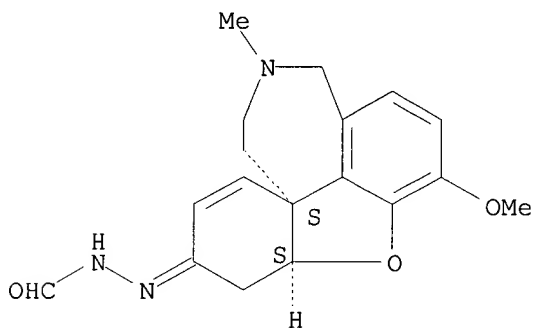
Relative stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hydrazinecarboxaldehyde,
 (4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-
 benzofuro[3a,3,2-ef][2]benzazepin-6-ylidene)-, (4aR*,8aR*)- (9CI)
 MF C18 H21 N3 O3

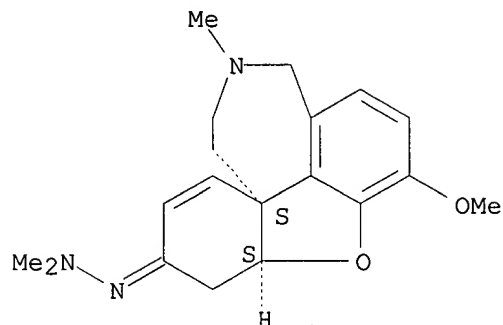
Relative stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
 methoxy-11-methyl-, dimethylhydrazone, [4aS-(4aR*,8aR*)]- (9CI)
 MF C19 H25 N3 O2

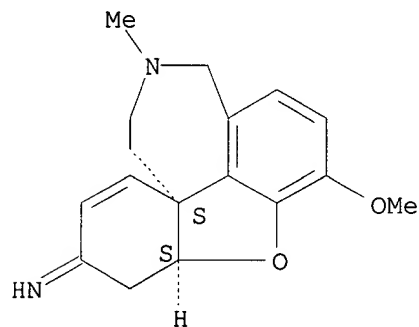
Relative stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-imine,
 4a,5,9,10,11,12-hexahydro-3-
 methoxy-11-methyl-, [4aS-(4aR*,8aR*)]- (9CI)
 MF C17 H20 N2 O2

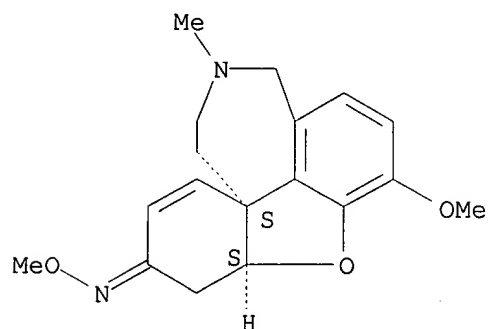
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
 methoxy-11-methyl-, O-methyloxime, [4aS-(4aR*,8aR*)]- (9CI)
 MF C18 H22 N2 O3

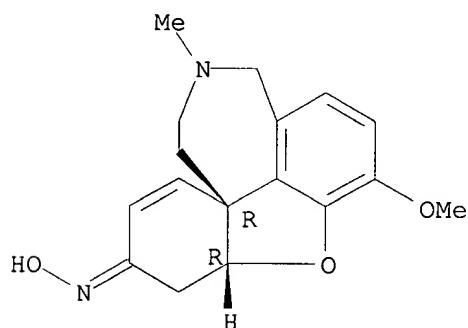
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, oxime, [4aR-(4aR*,8aR*)]- (9CI)
MF C17 H20 N2 O3

Absolute stereochemistry.
Double bond geometry unknown.

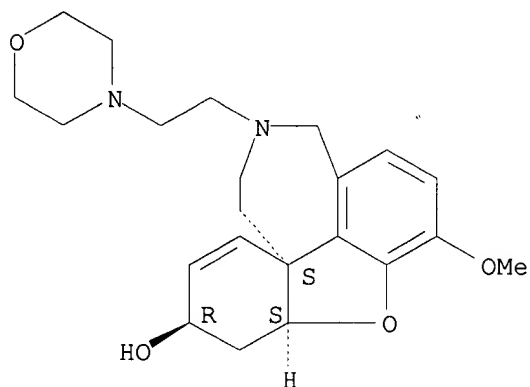


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-[2-(4-morpholinyl)ethyl]-, (4a.alpha.,6.beta.,8aR*)- (9CI)

MF C22 H30 N2 O4

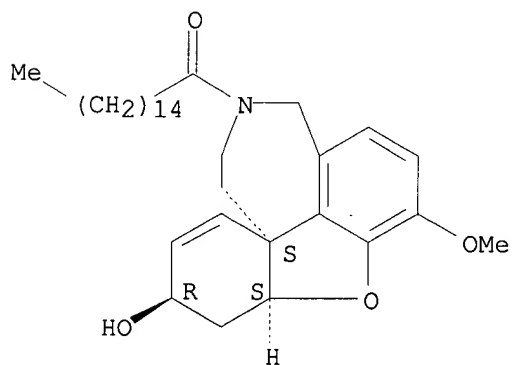
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-(1-oxohexadecyl)-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C32 H49 N O4

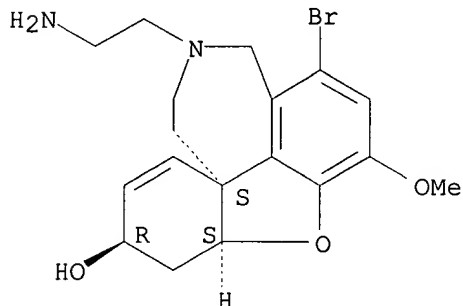
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-(2-aminoethyl)-1-bromo-
4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H23 Br N2 O3

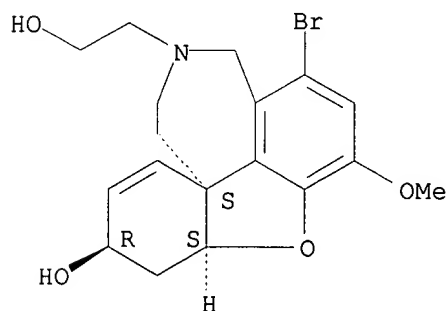
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-ethanol,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H22 Br N O4

Relative stereochemistry.

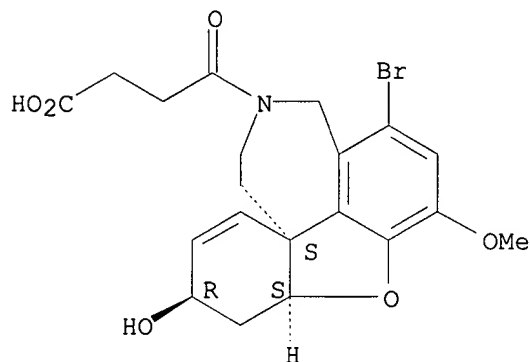


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-butanoic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.gamma.-oxo-,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H22 Br N O6

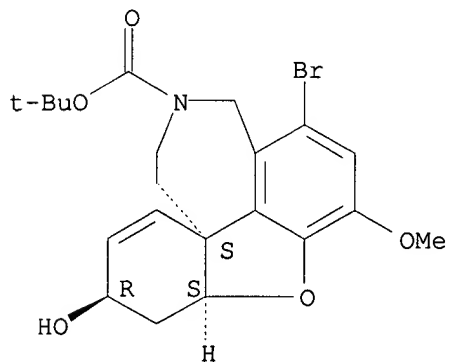
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxylic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, 1,1-dimethylethyl
ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H26 Br N O5

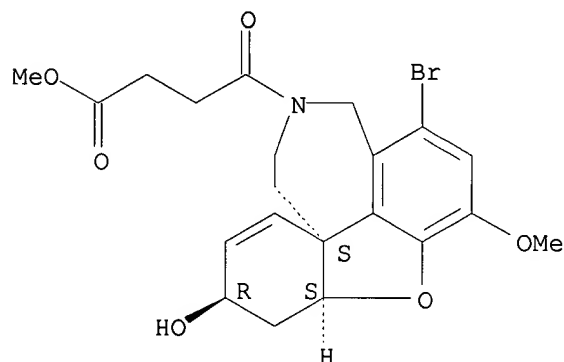
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-butanoic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-.gamma.-oxo-, methyl
ester, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C21 H24 Br N O6

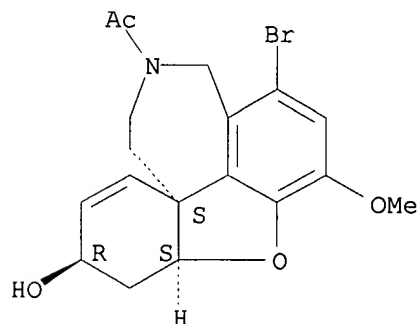
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 11-acetyl-1-bromo-
4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H20 Br N O4

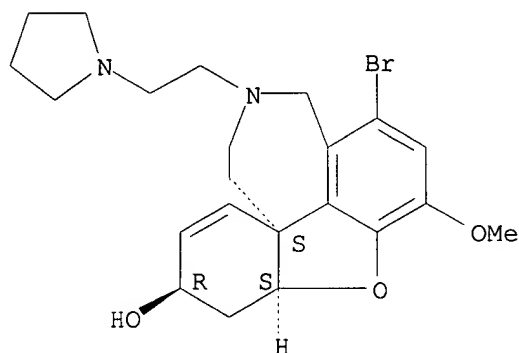
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-[2-(1-pyrrolidinyl)ethyl]-,
 (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C22 H29 Br N2 O3

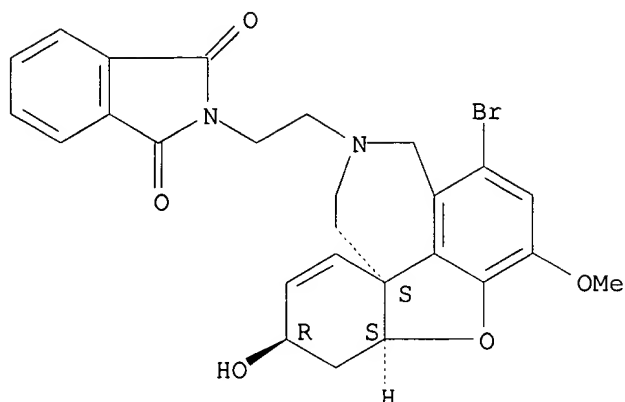
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1H-Isoindole-1,3(2H)-dione,
 2-[2-(1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-
 methoxy-6H-benzofuro[3a,3,2-ef][2]benzazepin-11(12H)-yl)ethyl]-,
 (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C26 H25 Br N2 O5

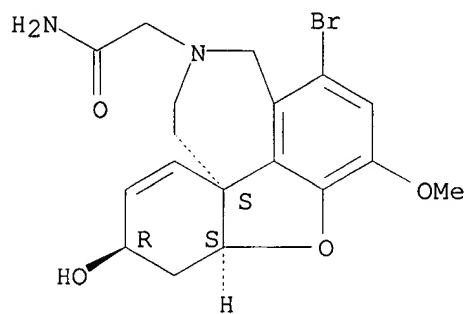
Relative stereochemistry. .



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetamide,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C18 H21 Br N2 O4

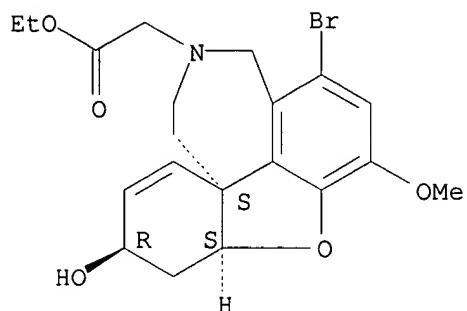
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-acetic acid,
1-bromo-4a,5,9,10-tetrahydro-6-hydroxy-3-methoxy-, ethyl ester,
(4a.alpha.,6.beta.,8aR*)- (9CI)
MF C20 H24 Br N O5

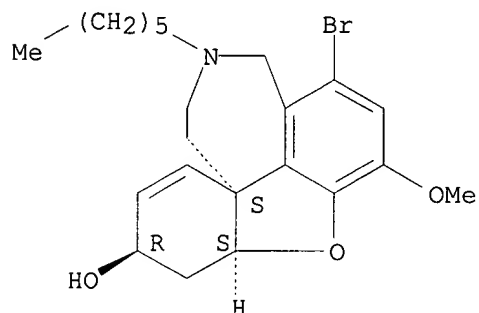
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-11-hexyl-
 4a,5,9,10,11,12-hexahydro-3-methoxy-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C22 H30 Br N O3

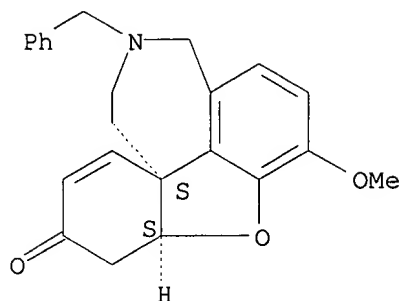
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-
 methoxy-11-(phenylmethyl)-, (4aR*,8aR*)- (9CI)
 MF C23 H23 N O3

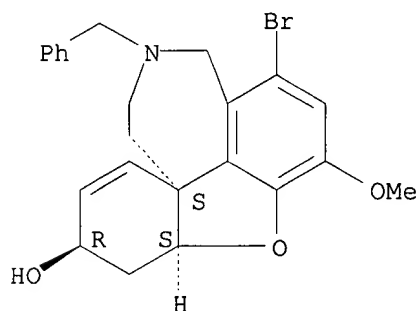
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-(phenylmethyl)-, (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C23 H24 Br N O3

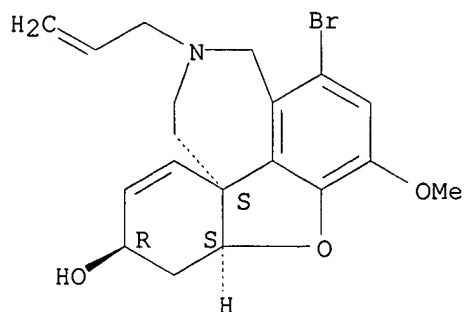
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-(2-propenyl)-, (4a.alpha.,6.beta.,8aR*)-(9CI)
 MF C19 H22 Br N O3

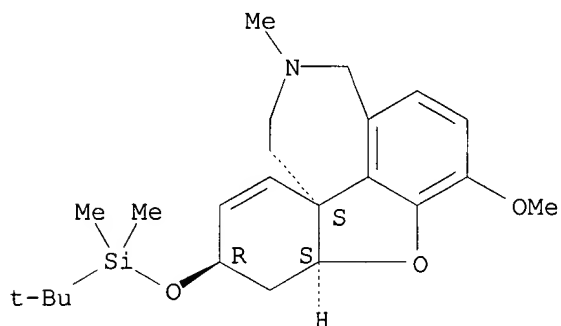
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine, 6-[[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C23 H35 N O3 Si

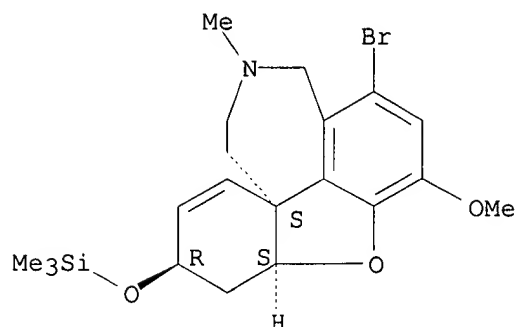
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine,
 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6-[(trimethylsilyl)oxy]-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C20 H28 Br N O3 Si

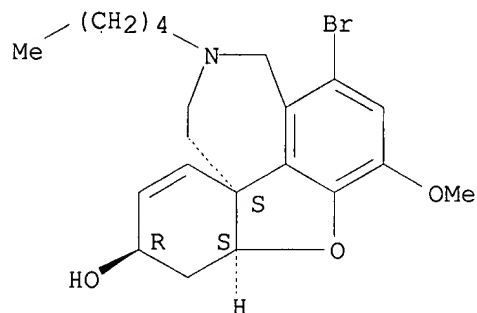
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-pentyl-, (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C21 H28 Br N O3

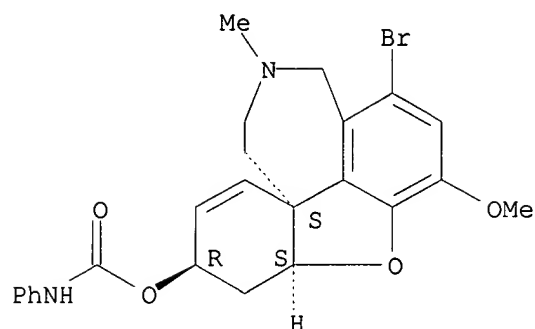
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-methyl-, phenylcarbamate (ester),
 (4a.alpha.,6.beta.,8aR*)- (9CI)
 MF C24 H25 Br N2 O4

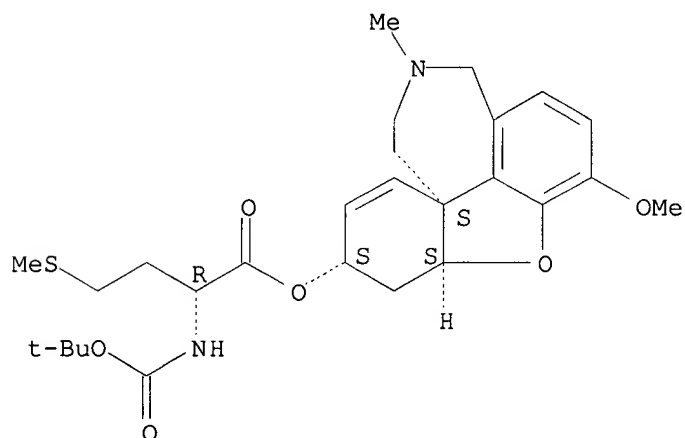
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN D-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-
 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
 ef][2]benzazepin-6-yl ester (9CI)
 MF C27 H38 N2 O6 S

Absolute stereochemistry.

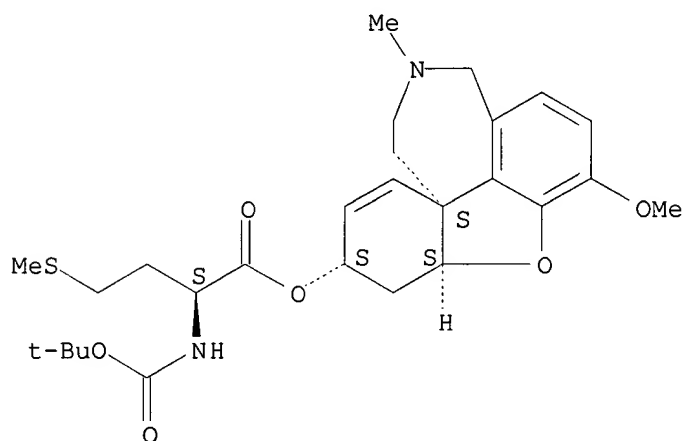


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN L-Methionine, N-[(1,1-dimethylethoxy)carbonyl]-, (4aS,6S,8aS)-

4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
ef][2]benzazepin-6-yl ester (9CI)
MF C27 H38 N2 O6 S

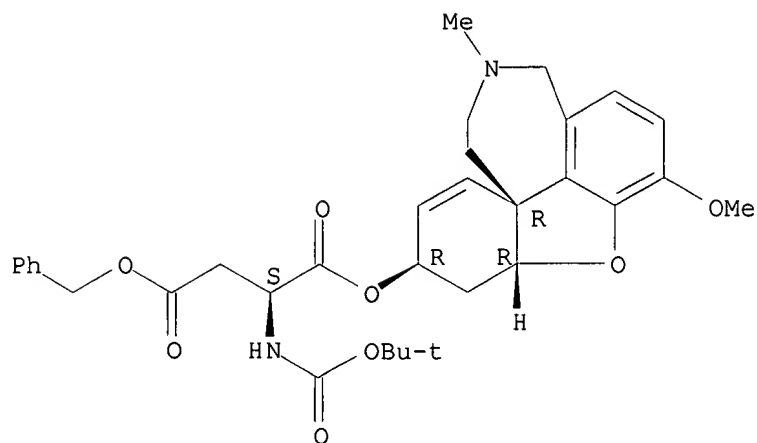
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN L-Aspartic acid, N-[(1,1-dimethylethoxy)carbonyl]-, 1-[(4aR,6R,8aR)-
4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-
ef][2]benzazepin-6-yl] 4-(phenylmethyl) ester (9CI)
MF C33 H40 N2 O8

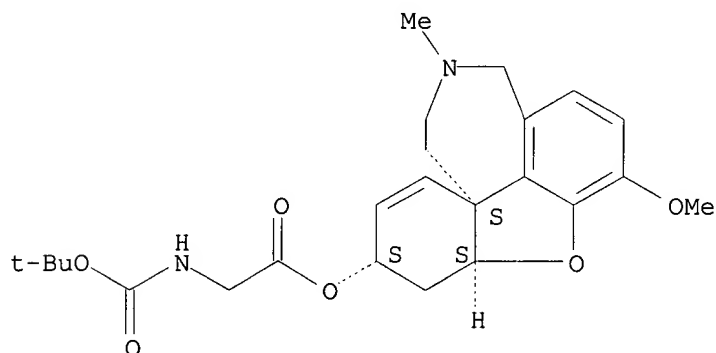
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Glycine, N-[(1,1-dimethylethoxy)carbonyl]-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester, [4aS-(4a.alpha.,6.alpha.,8aR*)]- (9CI)
 MF C24 H32 N2 O6

Absolute stereochemistry.



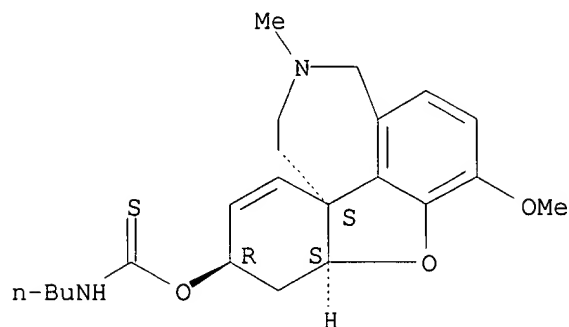
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Carbamothioic acid, butyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester,

Hong Liu

[4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C22 H30 N2 O3 S

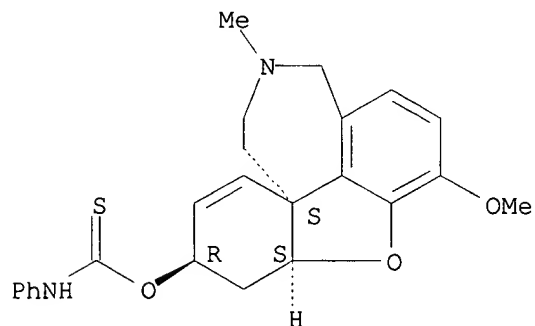
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamothioic acid, phenyl-, O-(4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl) ester,
[4aS-(4a.alpha.,6.beta.,8aR*)]- (9CI)
MF C24 H26 N2 O3 S

Absolute stereochemistry.

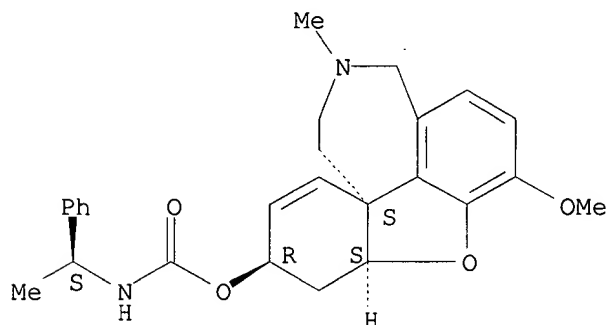


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester,
[4aS-[4a.alpha.,6.beta.(R*),8aR*]]- (9CI)
MF C26 H30 N2 O4

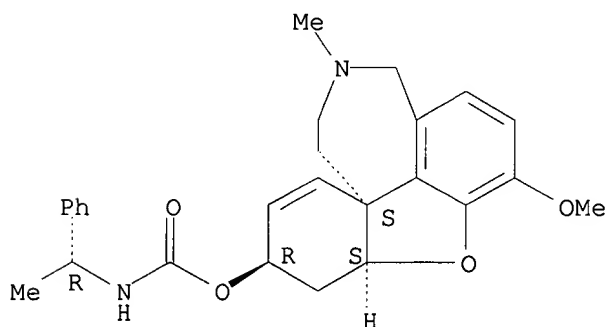
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, (1-phenylethyl)-, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-yl ester,
[4aS-[4a.alpha.,6.beta.(S*),8aR*]]- (9CI)
MF C26 H30 N2 O4

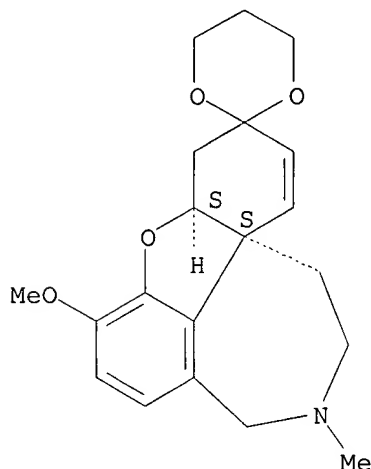
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxane],
 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR*,8aR*)- (9CI)
 MF C20 H25 N O4

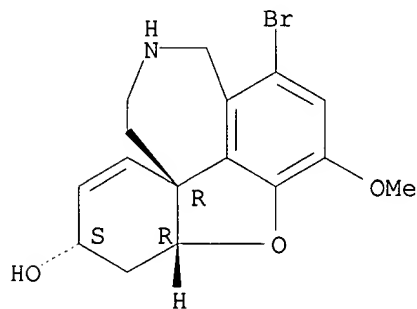
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-, [4aR-(4a.alpha.,6.beta.,8aR*)]- (9CI)
 MF C16 H18 Br N O3

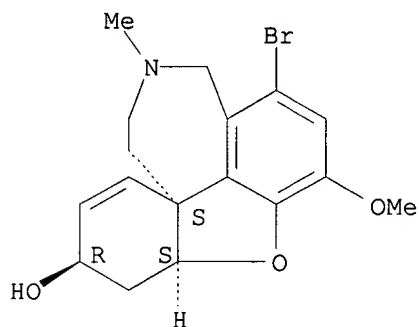
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI)
MF C17 H20 Br N O3

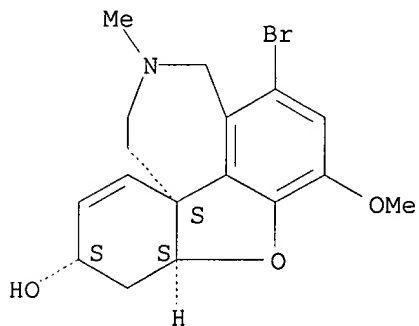
Absolute stereochemistry. Rotation (-).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
hexahydro-3-methoxy-11-methyl-, (4aR,6R,8aR)-rel- (9CI)
MF C17 H20 Br N O3

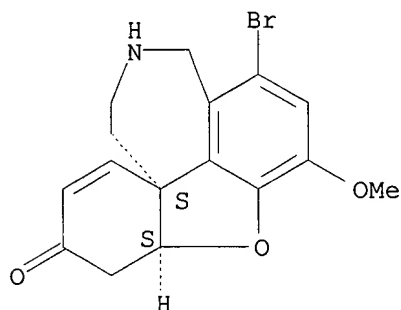
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-, (4aR,8aR)-rel- (9CI)
 MF C16 H16 Br N O3

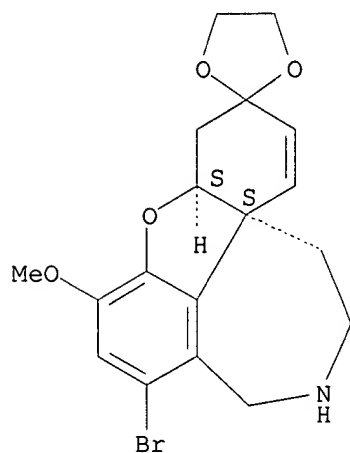
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane],
 1-bromo-4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR*,8aR*)- (9CI)
 MF C18 H20 Br N O4

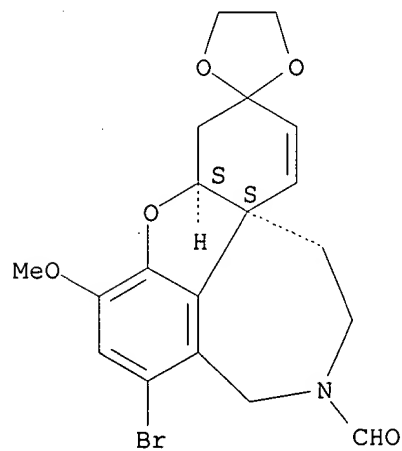
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Spiro[6H-benzofuro[3a,3,2-ef][2]benzazepine-6,2'-[1,3]dioxolane]-11(12H)-
 carboxaldehyde, 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-, (4aR*,8aR*)-
 (9CI)
 MF C19 H20 Br N O5

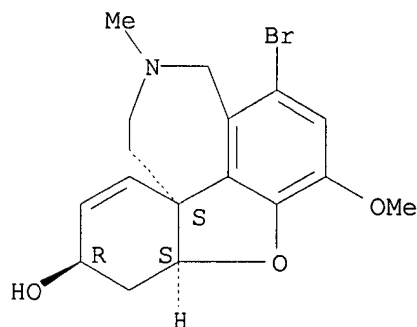
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 1-bromo-4a,5,9,10,11,12-
 hexahydro-3-methoxy-11-methyl-, (4aR,6S,8aR)-rel- (9CI)
 MF C17 H20 Br N O3

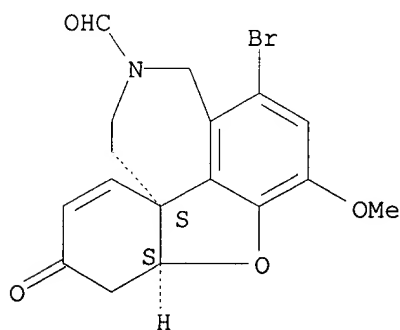
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

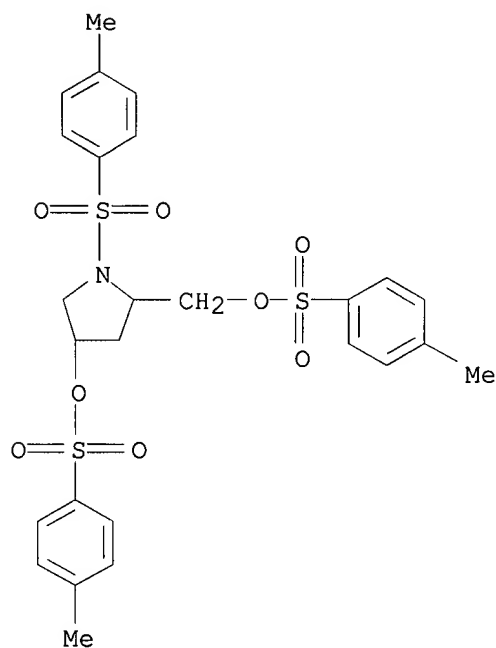
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepine-11(12H)-carboxaldehyde,
 1-bromo-4a,5,9,10-tetrahydro-3-methoxy-6-oxo-, (4aR,8aR)-rel- (9CI)
 MF C17 H16 Br N O4

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

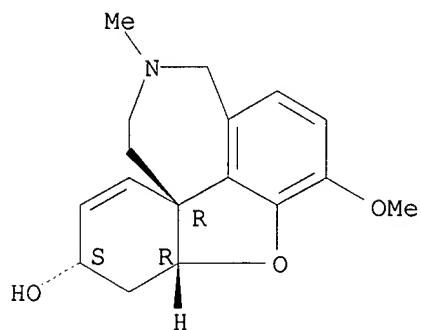
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 2-Pyrrolidinemethanol, 1-[(4-methylphenyl)sulfonyl]-4-[[4-methylphenyl)sulfonyl]oxy]-, 4-methylbenzenesulfonate (ester) (9CI)
 MF C26 H29 N O8 S3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aR,6S,8aR)- (9CI)
 MF C17 H21 N O3
 CI COM

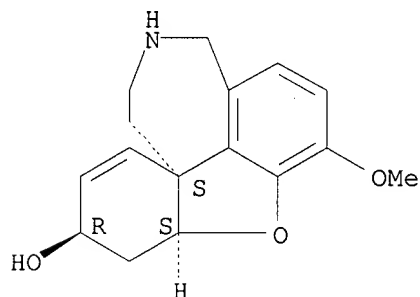
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-, (4aR,6S,8aR)-rel- (9CI)
 MF C16 H19 N O3

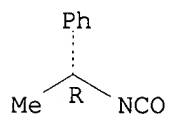
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzene, [(1R)-1-isocyanatoethyl]- (9CI)
 MF C9 H9 N O
 CI COM

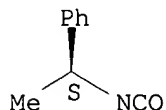
Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, [(1S)-1-isocyanatoethyl]- (9CI)
MF C9 H9 N O

Absolute stereochemistry. Rotation (-).



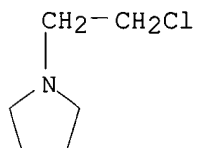
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazine, monohydrate (8CI, 9CI)
MF H4 N2 . H2 O
CI COM

H₂N-NH₂

● H₂O

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Pyrrolidine, 1-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
MF C6 H12 Cl N . Cl H

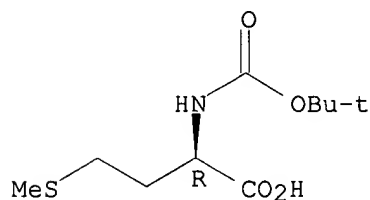


● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN D-Methionine, N-[(1,1-dimethylethoxy)carbonyl]- (9CI)
MF C10 H19 N O4 S
CI COM

Absolute stereochemistry.



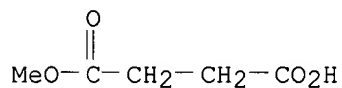
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Ethanamine, 2-chloro-N,N-dimethyl-, hydrochloride (9CI)
MF C4 H10 Cl N . Cl H

Me₂N-CH₂-CH₂Cl

● HCl

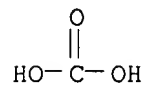
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Butanedioic acid, monomethyl ester (9CI)
MF C5 H8 O4
CI COM



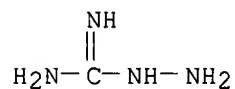
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbonic acid, compd. with hydrazinecarboximidamide (1:1) (9CI)
MF C H6 N4 . C H2 O3
CI COM

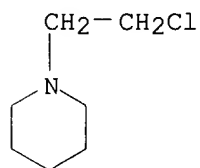
CM 1



CM 2

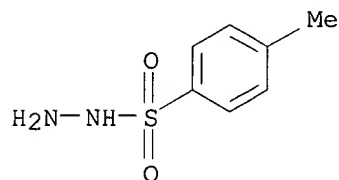


L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Piperidine, 1-(2-chloroethyl)-, hydrochloride (6CI, 7CI, 8CI, 9CI)
 MF C7 H14 Cl N . Cl H



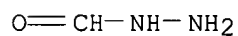
● HCl

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzenesulfonic acid, 4-methyl-, hydrazide (9CI)
 MF C7 H10 N2 O2 S
 CI COM



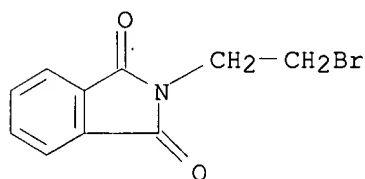
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazinecarboxaldehyde (9CI)
MF C H4 N2 O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

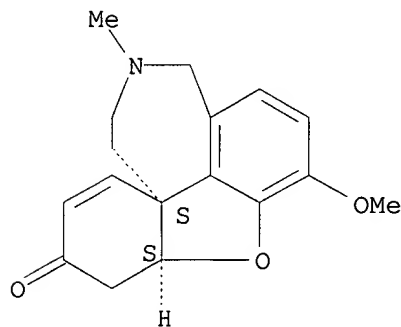
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 1H-Isoindole-1,3(2H)-dione, 2-(2-bromoethyl)- (9CI)
MF C10 H8 Br N O2

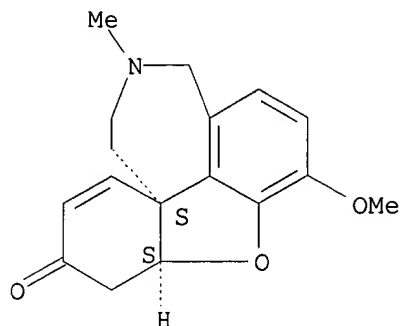


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-one, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,8aS)- (9CI)
MF C17 H19 N O3
CI COM

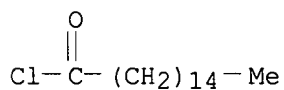
Absolute stereochemistry. Rotation (-).





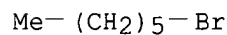
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hexadecanoyl chloride (9CI)
 MF C16 H31 Cl O



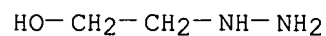
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Hexane, 1-bromo- (6CI, 8CI, 9CI)
 MF C6 H13 Br
 CI COM



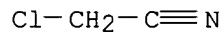
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Ethanol, 2-hydrazino- (6CI, 7CI, 8CI, 9CI)
 MF C2 H8 N2 O
 CI COM



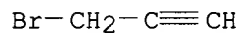
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Acetonitrile, chloro- (6CI, 8CI, 9CI)
 MF C2 H2 Cl N
 CI COM



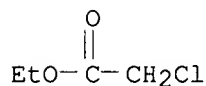
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN 1-Propyne, 3-bromo- (9CI)
 MF C3 H3 Br
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

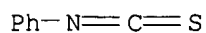
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Acetic acid, chloro-, ethyl ester (6CI, 8CI, 9CI)
 MF C4 H7 Cl O2
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

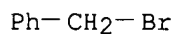
L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
 IN Benzene, isothiocyanato- (9CI)
 MF C7 H5 N S
 CI COM





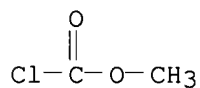
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Benzene, (bromomethyl)- (9CI)
MF C7 H7 Br
CI COM



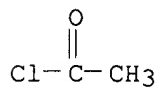
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Carbonochloridic acid, methyl ester (9CI)
MF C2 H3 Cl O2
CI COM



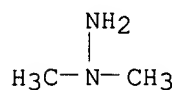
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Acetyl chloride (8CI, 9CI)
MF C2 H3 Cl O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 197 ANSWERS REGISTRY COPYRIGHT 2002 ACS
IN Hydrazine, 1,1-dimethyl- (8CI, 9CI)
MF C2 H8 N2
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> logoff

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE

ENTRY

0.76

TOTAL

SESSION

4.22

STN INTERNATIONAL LOGOFF AT 18:01:46 ON 04 OCT 2002